Phase I — Initial Site Investigation Report Former Nu-Style Property

87 Grove Street (Lots 22 & 27) Franklin, MA RTN 2-16694

Town of Franklin, Massachusetts

May 2008



317 Iron Horse Way, Suite 204 Providence, RI 02908



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1.0 GENERAL DISPOSAL SITE INFORMATION

1.1 Introduction

This report, prepared in accordance with the Massachusetts Contingency Plan (MCP) 310 CMR 40.0480, serves as documentation for a Phase I — Initial Site Investigation Report (Phase I Report) for response actions conducted at the disposal site (the "site") identified by the Massachusetts Department of Environmental Protection (MADEP) as Release Tracking Number (RTN) 2-16694. Fuss & O'Neill, Inc. (Fuss & O'Neill) prepared this report on behalf of the Town of Franklin, Massachusetts. Assessment activities were conducted as part of the Norfolk County Hazardous Materials and Petroleum Brownfield Assessment Programs funded under, two brownfield assessment grants from the United States Environmental Protection Agency (USEPA). Fuss & O'Neill's Limitations of Work Product are included as Appendix A.

1.2 <u>Site Description</u>

According to the Town of Franklin assessor's mapping, the subject property is comprised of approximately 1.20 acres on the following two parcels in the Town of Franklin, Norfolk County, Massachusetts:

<u>Table 1</u>
Summary of Parcels Comprising Subject Property

Plat/Lot	Address	Owner	Acreage*	Zoning
276/22	Grove Street	Town of Franklin	0.23	Business
276/27	Grove Street	Town of Franklin	0.97	Business

^{*} Acreage for the site was obtained from assessor's mapping.

The subject property's main address is 87 Grove Street and is located along the west side of Grove Street south of the intersection of Grove Street and Route 140. The location of the subject property is also defined as the following:

Latitude: 42°5′13.154″ North
Longitude: 71°25′39.790″ West

• UTM Coordinates (NAD 83 meters) = 4662290 North, 299210 East

A site location map depicting 500-foot and half-mile radii from the boundaries of the disposal site is attached as <u>Figure 1</u>. A site plan depicting the disposal site is presented as <u>Figure 2</u>.

A vacant, partially dilapidated two-story building with a footprint of approximately 11,800 square feet is situated on Lot 27, and a vacant one and one-half-story building with a footprint of approximately 4,000 square feet is located on Lot 22. A commercial development is currently proposed for the subject property. Formal development plans have not been established or approved by the Town of Franklin. An estimated 620 people



reside within a one-half mile radius of the subject property. The subject property is located in a mixed commercial and residential area in Franklin, Massachusetts.

No institutions are located within 500 feet of the site. The closest school, KinderCare Learning Center, is located more than 1,500 feet northeast of the subject property.

1.3 Natural Resource Areas

Mine Brook flows westward along the southern side of the Lot 27 building and turns northward to form the western boundary of Lot 22. Mine Brook flows generally northward to the Charles River.

According to the MADEP Bureau of Waste Site Cleanup Site Scoring Map, attached as Appendix B, no drinking water supplies are located within 500 feet of the disposal site.

No endangered species habitat, areas of critical environmental concern (ACEC) or certified vernal pools were identified located within 500 feet of the subject site. Mapped wetlands are located approximately 50 feet east and 400 feet northwest of the subject property, in addition to the wetlands associated with Mine Brook.

No Sole Source Aquifers, fish habitats, and habitats of Species of Special Concern or Threatened or Endangered Species were identified located within 500 feet of the disposal site. A local, state, and/or federal protected open space is located within approximately 50 feet of the southern boundary of the subject property.

1.4 Geologic and Physiographic Setting

The topography of the site is generally flat, except at the banks of Mine Brook, where the topography drops steeply three to five feet to the river bed (USGS, 1987). The regional topography is hilly and generally drains to Mine Brook.

Surficial material at the site was mapped as loamy udorthents, which generally consist of moderately coarse-grained, deep and moderately deep, fairly well-drained soils (USDA, 2006). Fill described as sand, gravel, silt, and, in some cases, wood and brick was observed to depths of up to 14 feet below grade during drilling conducted on the site as part of Site Investigation activities conducted by Fuss & O'Neill at the subject property.

Bedrock beneath the site was mapped as grayish-pink to greenish-gray, equigranular to slightly porphyritic, Dedham Granite (Zen, 1983). Bedrock was encountered at the site during drilling as part of Site Investigation activities at depths of between four and 12.5 feet below grade (fbg).

1.5 <u>Soil and Groundwater Categorization</u>

Site visitors may include both adults and children. Based on the foreseeable site use as a commercial development, the frequency of site use by adult workers is expected to be high, but with generally low intensity. High intensity use may be required of some workers in F:\P2005\0458\F10\Phase | SIR\Phase | SIR\Phase



landscaped areas. Therefore, in accordance with the MCP (310 CMR 40.0933), soil at paved areas of the site is categorized as S-3, and soil in unpaved portions of the site is categorized as S-2. However, since the future site use has not been definitively determined to date, we have conservatively compared soil analytical data at the site to S-1, S-2, and S-3 soil categories, as indicated in the reports attached as <u>Appendix C</u> and <u>Appendix D</u>.

In accordance with the MCP (310 CMR 40.0932), groundwater at the site was classified as GW-2/GW-3. All groundwater in the Commonwealth of Massachusetts is considered a potential source of discharge to surface water and shall be categorized, at a minimum, as category GW-3. Groundwater at the disposal site that is located within 30 feet of an occupied structure and is less than 15 fbg is additionally defined as GW-2. Although the existing buildings are not occupied, groundwater is conservatively being compared to GW-2 and GW-3 categories.

The subject property is not located within a MADEP Zone II (aquifer protection area), potentially productive aquifer, or other GW-1 inclusionary criteria; therefore, a classification of GW-1 does not apply to the property. The MADEP Bureau of Waste Site Cleanup Site Scoring Map is attached as <u>Appendix B</u>.

1.6 <u>Disposal Site Map</u>

A site plan depicting the subject property is provided as <u>Figure 2</u>. The site plan depicts the disposal site boundaries, boundaries of properties located within the disposal site, on-site buildings, floor and storm drains, subsurface utilities, the location of oil and/or hazardous materials (OHM) releases, and the location of monitoring wells, soil borings, and sediment and surface water samples installed or collected by Fuss & O'Neill.

2.0 DISPOSAL SITE HISTORY

2.1 Owner/Operator and Operations History

The site was acquired via tax title by the Town of Franklin as a result of foreclosure and tax title during 2002.

Historical topographic maps depict a building on both Lot 22 and Lot 27 by 1893. According to files available at the Town of Franklin offices, Unionville Woolen Mills operated on the subject property and on properties adjacent to the north, northeast, south, and southwest, likely since the site was first developed. Town property cards indicated that the current site buildings were originally constructed circa 1900 (Lot 27) and circa 1945 (Lot 22). Several additions appear to have been constructed onto both buildings.

A right-of-way currently located along the eastern boundary of Lot 22 is also known as "Old Grove Street." Grove Street was relocated from the right-of-way to the current location east of the subject property in the mid to late 1950s. At that time, the portion of the pond located on the subject property was partially filled and Mine Brook was relocated to flow to the south of the Lot 27 building, as shown in a 1968 Plan of Land prepared for Unionville Woolen Mills, Inc. The origin of the fill materials could not be determined during the Site F:\P2005\0458\F10\Phase I SIR\Phase I SIR\Phase I SIR\revised-040308-bek.doc



Investigation documented herein; however, we infer that the area was filled as part of the Grove Street relocation municipal project. The fill area is currently mostly paved for use by commercial businesses occupying Lot 26 (adjacent to the south) for parking.

Mapping available at the Town offices indicated that the Franklin Paint Company occupied the subject property and the parcel adjacent to the south at some point in the past, possibly in the 1950s. A 1956 plan prepared for the Franklin Paint Company depicted a dam on the south-central portion of Lot 27, at the eastern end of the reservoir. Until the early 1960s, the western end of Mine Brook Reservoir covered the eastern portion of Lot 27. The reservoir is referred to as a pond in subsequent mapping, which shows the pond partially beneath the Lot 27 building.

Carol and Richard Armstrong purchased the subject property in 1969 and used the property for jewelry manufacturing until the late 1980s under the names Nu-Style Company, Inc. and Image Jewelry. An elevated passageway (a covered pedestrian bridge) was constructed over Mill Brook circa 1969/1970. This bridge joined the Lot 27 building to the building located on Lot 26, adjacent to the south of the subject property. This bridge has since been demolished.

A 1975 plan for a proposed addition to the Lot 27 building indicated that this building was a manufacturing plant and the Lot 22 building was a garage. In 1978, the subject property was first listed in the Town Clerk's database for on-site storage of hazardous materials and USTs located on the property. The Lot 22 building was most recently used by a construction company for vehicular repair until it was vacated in 1989 (IES, 1990). Operations on both site parcels ceased in late 1989.

Aerial photographs from 1997 depicted the site buildings on Lots 22 and 27 and a covered pedestrian bridge over Mine Brook, connecting Lot 27 to the abutting parcel to the south (Lot 26). Vegetation was observed in the area immediately associated with Mine Brook in the 1997 aerial photograph. In 2001, Grove Street Towing and Tire operated out of the garage adjacent to the main site building. In the 2005 aerial photograph, the pedestrian bridge connecting Lot 27 and Lot 26 was not observed and the area on the south side of Mine Brook on Lot 27 was paved. Two additional truck-sized structures were present along the west and northeast sides of the building on Lot 27 in the 2005 aerial photograph. These structures were not observed during the Site Investigation activities documented herein.

A detailed record of ownership for the subject property was not readily available at the Town of Franklin Tax Assessor's office. It was determined from an outdated property card that Unionville Woolen Mills, Inc. owned the subject property from January 1962, or some date prior, until May 1969, upon which the subject property was sold to Richard and Carol Armstrong. Ownership of the site prior to 1962 was not readily determined.



2.2 Release History

2.2.1 UST Removal and CERCLA Removal Action

USTs were removed from the subject property in 1990, according to records maintained at the Town Clerk's office. The tanks included one 5,000-gallon, two 2,000-gallon, and one 1,000-gallon USTs.

USEPA conducted an inspection of the subject property on January 8, 1992. According to Town files, the inspection revealed the presence of full and partially full labeled drums and containers as well as drums and containers with undocumented material. The inspection also included the observation of seven process tanks in the former plating department which contained undocumented liquids and/or sludges. Some of the chemicals identified at the subject property included: sodium cyanide, chromic acid, potassium cyanide, perchloroethylene (tetrachloroethylene; PCE), zinc cyanide, nickel sulfate, and copper cyanide. Following the inspection, USEPA conducted Comprehensive Environmental Response, Compensation, and Liability Act (CERCLA) removal actions at the subject property in 1992. Removal actions included the removal and offsite disposal of hazardous waste, contaminated soil and debris, and product pumped from tanks located on the subject property.

2.2.2 IES, Inc. Site Investigation Reports

Portions of two reports prepared by IES, Inc. (IES) summarizing environmental investigations previously conducted on the subject property and on the parcel adjacent to the south were reviewed. The results of the investigations documented in these reports were summarized in a Phase II Environmental Site Assessment (ESA) prepared by Fuss & O'Neill in September 2007. A copy of the September 2007 Phase II ESA is attached as Appendix C. The conclusions of these reports include the following:

In January 1990, IES completed a report of a Chapter 21E Site Evaluation of 87 Grove Street for Home National Bank of Milford, Massachusetts. IES concluded that no releases of hazardous materials or petroleum products had occurred at the subject property; however, it is Fuss & O'Neill's opinion that the IES investigation was not adequate to definitively rule out releases on the subject property.

In July 1991, IES collected soil samples from four additional borings (B-1A through B-4A) to assess whether releases associated with USTs had occurred. Soil encountered at the site generally consisted of fill containing loam, sand, gravel, and, in some cases, brick and cinders. Fill materials were observed to depths of up to 8.5 feet below grade (fbg). Deeper soil consisted of very dense, fine-grained sand, silt, and gravel. Groundwater was encountered at depths of approximately 8.5 to 9 fbg. Monitoring wells were installed within the borings to allow for the collection of groundwater samples. IES identified releases of chlorinated solvents to soil and groundwater at boring location B-4A, which was advanced downgradient of USTs at the site and north of Mine Brook. Based on the apparent vertical distribution of VOCs in soil, IES inferred that the presence of VOCs was the result of a surface release.



2.2.3 Fuss & O'Neill Assessment Activities

Fuss & O'Neill conducted environmental assessment activities at the subject property in 2006 and 2007, under two brownfield assessment grants. The assessment activities were summarized in the following reports:

- Fuss & O'Neill, 2006. Phase I Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, May 2006.
- Fuss & O'Neill, 2007. Phase I Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, February 2007.
- Fuss & O'Neill, 2007. UST Closure Assessment Report, Former Nu-Style Company, Inc., Franklin, Massachusetts, July 2007. A copy of this report is included in <u>Appendix D</u>.
- Fuss & O'Neill, 2007. Phase II Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, September 2007. A copy of this report is included as <u>Appendix C</u>.
- Fuss & O'Neill, 2008. Phase II Environmental Site Assessment Addendum, Former Nu-Style Property, RTN 2-0016694, 87 Grove Street (Lots 22 and 27), Franklin, Massachusetts, February 2008. A copy of this report is included as <u>Appendix D</u>.

A Phase I ESA, prepared by Fuss & O'Neill in May 2006 and updated in February 2007, identified the following recognized environmental conditions (RECs) at the site:

- The site had a long history (at least 90 years) of manufacturing, including textiles and jewelry. Materials used and stored at the site associated with jewelry manufacturing included cyanides, metals, chlorinated solvents, and petroleum products. Additional substances associated with textile manufacturing were also likely used. Files indicated that numerous drums of hazardous waste and petroleum products were located outside of the site buildings.
- At least one UST was present on the western side of the Lot 27 building. In addition, a heating oil tank reportedly existed in an underground bunker on the same side of the building.
- A small tunnel containing slow-moving water was present beneath the Lot 22 building. A review of mapping on file at the Town Building Department suggested that the tunnel runs, or ran in the past, from Mine Brook and beneath the Lot 27 building to the Lot 22 building. There is the potential that the tunnel was used by the former woolen mill for direct waste disposal to Mine Brook prior to the realignment of the brook in the 1960s.



- Releases of chlorinated solvents to soil and groundwater were identified on Lot 26, which abutted the site to the south. This property was owned and occupied by the same entities that owned and operated the facilities at the site; therefore, there is the potential that similar releases have occurred at the site. Due to the proximity of this property to the site, there is the potential for releases that occurred on this property to adversely affect groundwater quality at the site.
- The southern portion of the site contained a pond that was filled circa 1960. The nature and origin of the fill were not known.

MCP Phase I Initial Site Investigation activities conducted by Fuss & O'Neill at the site are discussed in the Phase II ESA and Phase II ESA Addendum reports attached as <u>Appendix C and D</u>, respectively. These activities are briefly summarized in <u>Sections 3.0 and 4.0</u> below. Based on the initial environmental assessment results, a release notification form was submitted to MADEP on May 10, 2007. MADEP assigned release tracking number (RTN) 2-0016694 to the site.

2.3 OHM Use and Storage History

Historic manufacturing operations at the subject property required the use and storage of cyanides, metals, chlorinated solvents including trichloroethene (TCE) and PCE, chromic acid, nickel sulfate and petroleum products. Use of OHM ceased at the subject property in 1989.

According to Town records, four USTs and associated piping were removed from the subject property in 1990, reportedly including USTs containing chlorinated solvents. One 5,000 gallon, two 2,000 gallon and one 1,000 gallon USTs were reportedly removed. One heating-oil UST was removed from the site in May 2007, during the investigations conducted by Fuss & O'Neill and documented in the July 2007 UST Closure Assessment Report included in Appendix D.

Several documented releases have been identified at the site by Fuss & O'Neill and by previous investigators. No additional information was available regarding OHM usage, quantities, and storage locations, or regarding the age or volume of the former USTs reportedly removed in 1990.

2.4 Waste Management History

With the exception of the environmental media and hazardous waste managed during removal and site investigation activities described in <u>Section 2.2</u>, no information was available regarding the waste management history of the subject property.

2.5 <u>Environmental Permits and Compliance History</u>

The following permits were documented for the site:

MADEP Order of Conditions, File No. CE159-937, March 22, 2007.



A Tier II Compliance History (BWSC107B) transmittal form is included in Appendix E.

2.6 <u>Potentially Responsible Parties</u>

The Town of Franklin is the current owner of the site and is identified as a Potentially Responsible Party.

3.0 SUBSURFACE INVESTIGATION ACTIVITIES

Fuss & O'Neill conducted subsurface investigation activities on behalf of the County of Norfolk and Town of Franklin, to investigate the nature and extent of the compounds of concern in soil, groundwater, and sediment at the disposal site. Sampling locations are depicted on Figure 2. Site Investigation activities were conducted from November 2006 to May 2007 in accordance with USEPA-approved Site-Specific Quality Assurance Project Plans (QAPPs). These activities were summarized in the September 2007 Phase II ESA included as Appendix C and the February 2008 Phase II ESA Addendum included as Appendix D. Please refer to these reports for details regarding the scope of work, summary of results, and data analysis of the subsurface investigation activities conducted at the disposal site. These reports are also briefly summarized herein.

Investigation and sampling activities conducted by Fuss & O'Neill included the following:

- A ground-penetrating radar (GPR) survey over portions of the subject property to evaluate for the potential presence of suspected USTs.
- Advancement of a total of 15 shallow soil borings and two deep soil borings.
- A soil sampling program that included the collection of a total of 29 soil samples, including quality control samples, from the soil borings. Soil boring logs depicting sample recovery amounts, material descriptions, graphic logs, soil codes, and photoionization detector (PID) soil screening results were included in the attached reports. A summary of soil sampling activities, including the requested analytical parameters, is included in <u>Table 2</u> below:

<u>Table 2</u> Summary of Soil Sampling Activities

Date	Sampling Location	Sample ID	Sample Depth (fbg)	Analyses
	B-02	30-03; 30-04	0-2; 5-7	VOCs, PP13 metals plus barium,
	B-04	30-07; 30-08	0-2; 5-6	total cyanide, PCBs, VPH, EPH
	B-05	30-09; 30-10	0.4-2; 5-7	
11/30/2006	B-06	30-11; 30-12	0-0.5; 1-2	
11/30/2000	MW-1	30-01; 30-02	0.5-2; 3-5	
	MW-2	30-05; 30-06	0-3; 5-7	
	MW-3	30-13; 30-14	0-2; 5-7	
	MW-4	30-15; 30-16*; 30-17	0-2; 0-2; 5-7	
12/1/2006	B-10	01-21; 01-22	0-2; 5-7	
	B-11	01-23; 01-24	0-2; 5-7	



Date	Sampling Location	Sample ID	Sample Depth (fbg)	Analyses
	B-12	01-25	0-2	
	MW-5	01-19; 01-20	0.4-2; 5-7	
10/31/2007	MW-13	31-01	10-12	VOCs
	B-15	01-03	2-4	VOCs
11/1/2007	MW-17	01-04; 01-05; 01-06*	0.3-2; 6-8; 6-8	VOCs, PP13 metals plus barium, VPH, EPH

Notes: B = soil boring

MW = monitoring well

Sample ID = Only the last 4 digits of the sample number are listed.

* indicates duplicate sample

VOCs = volatile organic compounds
PP13 metals = 13 Priority Pollutant metals
PCBs = polychlorinated biphenyls
VPH = volatile petroleum hydrocarbons
EPH = extractable petroleum hydrocarbons

- Installation of seven shallow overburden monitoring wells and two bedrock monitoring wells.
- Collection of groundwater samples from the monitoring wells. A summary of groundwater sampling activities, including the requested analytical parameters, is included in Table 3 below:

<u>Table 3</u> Summary of Groundwater Sampling Activities

Date	WellID	Sample ID	Analyses
	MW-1	08-27; 08-28*	
	MW-2	08-30	
12/8/06	MW-3	08-32	VOCs, PP13 metals plus barium, VPH, EPH
	MW-4	08-29	
	MW-5	08-31	
	MW-1	06-03	VOCs, total RCRA 8 metals, VPH, EPH
	MW-2	07-10	
	MW-3	07-11	VOCs, dissolved RCRA 8 metals
11/6/2007-	MW-5	07-09	
11/7/2007	MW-13	06-05; 06-06*	
	MW-14	06-04	VOCs, total RCRA 8 metals
	MW-16	07-08	
	MW-17	06-01; 06-02*	VOCs, total RCRA 8 metals, VPH, EPH

Notes: MW = monitoring well

Only the last four digits of the sample number are shown.

* indicates duplicate sample

VOCs = volatile organic compounds PP13 metals = 13 Priority Pollutant metals VPH = volatile petroleum hydrocarbons EPH = extractable petroleum hydrocarbons

RCRA 8 metals = Eight Resource Conservation and Recovery Act metals

 Collection of nine sediment samples, including quality control samples, from the banks of Mine Brook. A summary of sediment sampling activities, including the requested analytical parameters, is included in <u>Table 4</u> below:



<u>Table 4</u>
Summary of Sediment Sampling Activities

Date	Location	Sample ID	Analyses
	SD-1	SD-1	
4/26/2007	SD-2	SD-2	VOCs, PP13 metals plus barium, VPH, EPH, PCBs,
47 207 2007	SD-3*	SD-3*	cyanide
	SD-4	SD-4	
	SD-5	SD-5	
10/25/2007	SD-6	SD-6	SVOCs
	SD-7	SD-7	

Notes: SD = sediment sample

Only the last four digits of the sample number are shown.

* indicates duplicate sample

VOCs = volatile organic compounds PP13 metals = 13 Priority Pollutant metals VPH = volatile petroleum hydrocarbons

EPH = extractable petroleum hydrocarbons

PCBs = polychlorinated biphenyls

SVOCs = semi-volatile organic compounds

 Collection of five surface water samples, including quality control samples, from Mine Brook. A summary of surface water sampling activities, including the requested analytical parameters, is included in <u>Table 5</u> below:

<u>Table 5</u> Summary of Surface Water Sampling Activities

Date	Location	Sample ID	Analyses	
4/26/2007	SW-1	SW-1		
	SW-2	SW-2	VOCs, PP13 metals plus barium, VPH, EPH	
	SW-3*	SW-3*	VOC3, 11 13 metals plus bandin, V111, E111	
	SW-4	SW-4		

Notes: SW = surface water sample

Only the last four digits of the sample number are shown.

* indicates duplicate sample

VOCs = volatile organic compounds
PP13 metals = 13 Priority Pollutant metals
VPH = volatile petroleum hydrocarbons
EPH = extractable petroleum hydrocarbons

Closure of a 5,000-gallon #2 heating oil underground storage tank (UST) on May 1 and 2, 2007. Closure activities were completed by TMC Services, Inc. (TMC) of Bellingham, Massachusetts and observed by Fuss & O'Neill. Six confirmation samples were collected from the limits of the excavation for analysis of PP13 metals VOCs, and petroleum hydrocarbons (MADEP EPH and VPH Methods with target compounds). A UST Closure Assessment Report (CAR) was prepared by Fuss & O'Neill and submitted to MADEP in July 2007. A copy of the UST CAR was included in the September 2007 Phase II ESA attached as Appendix C.



4.0 SUBSURFACE INVESTIGATION RESULTS

A summary of subsurface investigation activities conducted at the site by Fuss & O'Neill is included in <u>Section 3.0</u>. Soil boring logs, monitoring well completion logs, and laboratory analytical reports were included in the respective reports included as <u>Appendix C</u> and <u>Appendix D</u>.

4.1 Soil Characterization

In general, the soil within soil borings advanced at the site was observed to consist of mainly fine to medium sand, with varying proportions of gravel and silt. Apparent fill material containing metal slag and coal and/or coal ash was observed in soil borings advanced north of Mine Brook, and was concentrated in the upper two feet of soil. Soil boring MW-17 ended in a soil horizon consisting predominately of silt and clay from 12 to 14 feet below grade. The silt and clay horizon may represent the sediment/water interface of the former pond that existed at this location prior to the historical placement of fill material.

Two soil horizons consisting predominately of silt were encountered in soil boring MW-05. The upper silty horizon had a minimum thickness of four feet, and occurred in the 5-10 foot depth interval. The lower silty horizon occurred in the 10-12 foot depth interval, had a minimum thickness of 0.8 feet, and may have extended beyond the maximum boring depth.

Bedrock encountered at the site was described as granite and was observed as shallow as four fbg at soil boring B-15.

4.2 <u>Soil Analytical Results</u>

Twenty-nine soil samples were collected from 16 soil borings throughout the site. The samples were collected from surficial soil, from vadose zone soil directly above the water table, or from soil horizons exhibiting the highest PID readings. Laboratory analytical results of soil samples collected from on-site soil borings documented the presence of the following analytes in soil at concentrations above laboratory reporting limits:

<u>Table 7</u>
Summary of Detected Compounds in Soil Samples

Metals (Method 6010)	EPH/PAH (MADEP EPH Method)	VPH (MADEP VPH Method)
Antimony	C19 to C36 Aliphatic Hydrocarbons	C9 to C12 Aliphatic Hydrocarbons
Arsenic	C11 to C22 Aromatic Hydrocarbons	
Barium	2-Methylnaphthalene	VOCs (Method 8260B)
Beryllium	Acenaphthene	Acetone
Cadmium	Acenaphthylene	M/P-xylenes
Chromium	Anthracene	Methyl Ethyl Ketone
Copper	Benzo(a)anthracene	Naphthalene
Lead	Benzo(a)pyrene	Tetrachloroethene (PCE)
Nickel	Benzo(b)fluoranthene	Toluene
Selenium	Benzo(k)fluoranthene	Trichloroethene (TCE)
Silver	Chrysene	
Thallium	Fluoranthene	



Metals (Method 6010)	EPH/PAH (MADEP EPH Method)	VPH (MADEP VPH Method)
Zinc	Fluorene	
Mercury (Method 7471)	Indeno(1,2,3-cd)pyrene	
	Phenanthrene	
	Pyrene	

Notes: PAH = polycyclic aromatic hydrocarbons

4.3 <u>Groundwater Analytical Results</u>

Sixteen groundwater samples were collected from eight monitoring wells at the site. Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of the following analytes in groundwater at concentrations above laboratory reporting limits:

<u>Table 8</u>
Summary of Detected Compounds in Groundwater Samples

Metals (Method 6010)	VOCs (Method 8260)
Barium, Total & Dissolved	Methyl tert butyl ether (MTBE)
Beryllium, Total	cis-1,2-Dichloroethene
Cadmium, Total	1,1,1-trichloroethane
Chromium, Total	PCE
Copper, Total	TCE
Lead, Total & Dissolved	
Nickel, Total	
Zinc, Total	

4.4 <u>Depth to Groundwater and Groundwater Flow</u>

Depth to groundwater in the on-site monitoring wells was measured by Fuss & O'Neill on several occasions. Based on the depth to water measurements collected from the monitoring wells and the site survey data, the direction of shallow groundwater flow at the site was generally to the south-southwest, as depicted on equipotential contour maps of shallow groundwater presented as <u>Figure 3</u> in both the Phase II ESA (<u>Appendix C</u>) and the Phase II ESA Addendum (Appendix D). Both figures are also included herein.

4.5 <u>Sediment Analytical Results</u>

Nine sediment samples were collected from seven locations along Mine Brook. Laboratory analytical results of sediment samples collected from Mine Brook documented the presence of the following analytes in sediment at concentrations above laboratory reporting limits:

<u>Table 10</u> Summary of Detected Compounds in Sediment Samples

Metals (Method 6010)	EPH/PAH (MADEP EPH Method)	VOCs (Method 8260B)
Arsenic	Acenaphthylene	Acetone
Barium	Anthracene	PCE
Beryllium	Benzo(a)anthracene	TCE
Cadmium	Benzo(a)pyrene	



Metals (Method 6010)	EPH/PAH (MADEP EPH Method)	VOCs (Method 8260B)
Chromium	Benzo(b)fluoranthene	
Copper	Benzo(k)fluoranthene	
Lead	Chrysene	
Nickel	Dibenzo(a,h)anthracene	
Thallium	Fluoranthene	
Zinc	Fluorene	
	Indeno(1,2,3-cd)pyrene	
	Phenanthrene	
	Pyrene	

4.6 <u>Surface Water Analytical Results</u>

Five surface water samples were collected from four locations along Mine Brook. Laboratory analytical results of surface water samples collected from Mine Brook documented the presence of the following analytes in surface water at concentrations above laboratory reporting limits:

<u>Table 12</u> Summary of Detected Compounds in Surface Water Samples

Metals (Method 6010)	VPH (MADEP VPH Method)
Barium	MTBE
Copper	
Lead	
Zinc	

4.7 <u>UST Confirmatory Soil Sample Results</u>

Six confirmation soil samples were collected from the limits of the UST grave following tank removal. Laboratory analytical results of the confirmatory soil samples collected from the UST grave documented the presence of the following analytes in soil at concentrations above laboratory reporting limits:

Table 14
Summary of Detected Compounds in Confirmation Soil Samples

Metals (Method 6010)	EPH/PAH (MADEP EPH Method)	VOCs (Method 8260B)
Arsenic	C9-C18 Aliphatic Hydrocarbons	Acetone
Barium	C19-C36 Aliphatic Hydrocarbons	PCE
Beryllium	C11-C22 Aromatic Hydrocarbons	TCE
Cadmium	Acenaphthylene	1,1,1-trichloroethane
Chromium	Anthracene	
Copper	Benzo(a)anthracene	
Lead	Benzo(a)pyrene	
Nickel	Benzo(b)fluoranthene	
Zinc	Benzo(k)fluoranthene	
	Fluoranthene	
	Fluorene	
	Indeno(1,2,3-cd)pyrene	
	Phenanthrene	
	Pyrene	



5.0 EVALUATION OF ANALYTICAL RESULTS

5.1 Data Verification

Procedures and methodologies for the collection and analyses of soil, groundwater, sediment and surface water samples were performed consistent with the MCP (310 CMR 40.0017). Analytical data were developed and reviewed in accordance with MADEP's Compendium of Quality Assurance and Quality Control Requirements and Performance Standards for Selected Analytical Methods (the CAM).

Presumptive Certainty was obtained for each data set collected as part of the subsurface investigation. Documentation was provided by Premier along with narrative summaries in the respective reports attached as <u>Appendix C</u> and <u>Appendix D</u>.

5.2 Soil

In accordance with 310 CMR 40.0361, the MCP Method 1 S-1, S-2, and S-3 standards for GW-2 and GW-3 areas were applied to soil samples obtained on the site. A summary of soil analytical results for samples collected by Fuss & O'Neill was included in each report included as <u>Appendix C</u> and <u>Appendix D</u>.

In summary, the concentrations of PCE and TCE exceeded the applicable standards in soil samples collected from borings B-04, B-06, and B-15. Additionally, the concentration of TCE in the soil sample collected from boring B-10 exceeded applicable standards. These soil samples were collected from the vadose zone to as deep as six fbg. These results suggest that these compounds were introduced to the subsurface via surficial releases.

The concentration of lead in soil samples collected from boring MW-05 up to seven fbg exceeded the applicable criteria. The concentration of beryllium in the soil sample collected from boring B-10 exceeded the S-2/GW-2 and S-2/GW-3 standards.

5.3 Groundwater

In accordance with 310 CMR 40.0362, the MCP Method 1 Groundwater standards for GW-2 and GW-3 areas were applied to groundwater samples obtained on the site. Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of lead at levels in excess of the GW-3 criteria in samples collected from monitoring wells MW-1, MW-3, and MW-5. The VOC compounds PCE and TCE were reported at levels in excess of the GW-2 criteria in samples collected from monitoring wells MW-3, MW-4, and MW-13. Additionally, the concentration of TCE in the sample collected from monitoring well MW-16 exceeded the GW-2 standard.

5.4 Sediment

Sediment analytical results were compared to the sediment screening values. Laboratory analytical results of sediment samples collected from Mine Brook documented the presence of several EPH or SVOC compounds in samples SD-01, SD-05, SD-06, and SD-07 at levels F:\P2005\0458\F10\Phase I SIR\Phase I SIR\revised-040308-bek.doc



in excess of the sediment screening values. The detected compounds are collectively identified as PAHs. These sediment samples were collected in close proximity to each other and were collected from the western portion of the site in the vicinity of sediment sample SD-01.

5.5 Surface Water

Surface water analytical results were compared to the MCP Method 1 Groundwater standards for GW-3 areas as well as the USEPA Chronic Criteria Continuous Concentrations, as documented in Table 10 of the February 2008 Phase II ESA Addendum attached as <u>Appendix D</u>. Laboratory analytical results of surface water samples collected from Mine Brook did not document the presence of compounds at levels in excess of these criteria.

5.6 <u>UST Confirmatory Soil Samples</u>

In accordance with 310 CMR 40.0361, the RCS-1 reporting category and the MCP Method 1 standards for S-1, S-2, and S-3 for GW-1, GW-2, and GW-3 areas were applied to confirmatory soil samples obtained on the subject site associated with the UST closure.

Laboratory analytical results of soil samples collected from the limits of the tank grave did not document the presence of compounds at levels in excess of the applicable criteria.

6.0 CONCEPTUAL SITE MODEL

A conceptual site model (CSM) has been developed for the disposal site based on the nature and source of the release, geologic and hydrogeologic conditions, historical site uses, current uses and foreseeable site uses. Available site data, including data presented herein and generated in previous investigations by Fuss & O'Neill, was evaluated in developing the CSM. The CSM was used to develop conclusions regarding the apparent extent of contamination, the media affected by the releases, and sufficiency of investigations. A discussion of the source, site hydrogeology, migration pathways, and the nature and extent of contamination follows.

6.1 <u>Site Hydrogeological Characteristics</u>

The disposal site was underlain by fine to medium sand, with varying proportions of gravel and silt. Apparent fill material containing metal slag and coal and/or coal ash was observed in soil borings advanced north of Mine Brook, and was concentrated in the upper two feet of soil. The soil deposits extended to at least 14 fbg based on the deepest soil boring (MW-17). No wide-ranging impermeable confining layer was identified in the available soil data at the site. Bedrock composed of granite was encountered as shallow as four fbg at soil boring B-15.

Two groundwater sampling events have been conducted at the disposal site and provide data related to seasonal groundwater occurrence and flow. The depth to groundwater ranged from approximately 4.2 fbg to approximately 9.2 fbg across the site. Shallow groundwater F:\P2005\0458\F10\Phase I SIR\Phase I SIR\revised-040308-bek.doc



flow direction was generally to the south-southwest toward Mine Brook. An equipotential contour map of shallow groundwater was presented as <u>Figure 3</u> in both the Phase II ESA and Phase II ESA Addendum. Both figures are also attached herein.

Two monitoring wells, designated MW-13, and MW-14, were advanced and screened in bedrock at the first water-bearing fracture encountered during drilling. Based on the groundwater elevations in these wells, groundwater flow direction in the bedrock aquifer was generally to the southward in the direction of Mine Brook. Groundwater elevations at these bedrock wells compared to the elevation of shallow groundwater in overburden monitoring wells indicated an upward vertical hydraulic gradient from the bedrock aquifer to the overburden aquifer. However, the hydraulic connection of the bedrock and overburden aquifers was not established during this investigation.

Based on a review of MCP GW-1 inclusionary criteria and the Site Scoring Map, attached as <u>Appendix B</u>, groundwater at the site and the surrounding area was not used for drinking water and was not considered a potential drinking water source.

6.2 <u>Contaminant Sources</u>

Based on the investigations and reports summarized herein, the following releases were identified:

- Soil: chlorinated VOCs (PCE and TCE), beryllium, and lead
 - The highest concentrations of VOCs were identified in soil samples collected proximal to the western and northern edges of the Lot 27 building at two general locations:
 - § In shallow soil in the vicinity of B-06, B-10, and B-15. The source of VOCs in this area may be attributed to historic surficial spills that impacted shallow soil.
 - § In shallow soil and soil at the approximate depth of the water table at soil boring B-04. The source of VOCs in this area may be attributed to both historic surficial spills as well as a potential additional source of VOCs in groundwater at depth.
 - The source of beryllium in the soil sample collected from boring B-10 and lead in the soil sample collected from boring MW-05 is likely attributed to natural sources and/or historic filling practices at the site.
- Shallow Overburden Groundwater: chlorinated VOCs (PCE and TCE) and lead
 - VOCs in shallow groundwater were identified in samples collected from monitoring wells MW-03, MW-04, and MW-16. These wells were located in the vicinity of soil containing VOCs, likely as a result of surficial spills that



- migrated to groundwater. The dissolved-phase VOCs in these areas indicated that VOCs migrated to groundwater as a result of the historic spills.
- Lead in shallow groundwater was identified in samples collected from monitoring wells MW-01, MW-03, and MW-05. The source of lead in shallow groundwater was attributed to leaching of lead to groundwater from fill material deposited at the site.
- Bedrock Groundwater: chlorinated VOCs (PCE and TCE)
 - OVOCs in bedrock groundwater were identified in the sample collected from monitoring well MW-13. The source of VOCs in this monitoring well may be attributed to a combination of both surficial releases of VOCs to the subsurface as well as a potential additional source from within the on-site building. The additional source has not been confirmed, and additional investigation may be warranted following the demolition of the building.
- Sediment: PAHs
 - PAHs in sediment were identified in the vicinity of sediment samples SD-01, SD-05, SD-06, and SD-07 at the western portion of the site and at downstream locations of Mine Brook.
 - The source of PAHs in sediment was attributed to historic filling practices at the site in which Mine Brook was channelized/re-located and former surface waters at the site were filled in; or to historic discharges from the former manufacturing facility.
- Surface Water: No releases to surface water were identified.
- UST Confirmatory Soil Samples: No releases to environmental media were identified associated with the former UST located at the southwest portion of the building.

The use and storage of chlorinated solvents at the site was discontinued when the site was vacated in 1989, thus eliminating the principal surface source of VOC contamination to site soil and groundwater.

6.3 Nature and Extent of Contamination

Based on the results of soil, groundwater, sediment, and surface water monitoring conducted at the site the lateral extent of the disposal site has generally been delineated and is shown on <u>Figure 2</u>. The nature and extent of contamination at the disposal site, based on the results of all Site Investigation activities conducted at the disposal site by Fuss & O'Neill, is detailed in <u>Section 4.0</u> and <u>Section 5.0</u>. However, additional delineation of the source and/or extent of VOCs in bedrock groundwater as well as PAHs in sediment may be warranted.



6.4 <u>Migration Pathways and Exposure Potential</u>

A description of the migration pathways associated with each of the types of environmental media affected by the releases at the site, sediment, soil, and groundwater, are summarized in the following paragraphs. These potential pathways are the primary methods for migration of site-related contaminants of concern (COC).

Migration and mobilization of COC in shallow soil may occur via infiltration of stormwater though vadose zone soil containing COC and via migration of shallow groundwater through saturated soil containing COC. VOCs in shallow soil can also volatilize and migrate into building structures located above soil containing VOCs.

Migration of COC in shallow overburden groundwater may occur via horizontal migration though saturated overburden toward Mine Brook. Although the data presented in <u>Section 6.1</u> indicates that the vertical gradient in the central portion of the site is generally upward, the increasing concentration of dissolved VOCs with depth indicates a chlorinated VOC pathway downward to the deeper bedrock aquifer. The relative concentrations of COC in shallow and deep groundwater are indicative of a potential unidentified source within or beneath the abandoned building and a potentially historical direct release from the facility to the bedrock aquifer. Migration of COC in the deeper bedrock aquifer is considered to be via fracture flow, and may not be well defined by the equipotential contours representing shallow groundwater gradients. VOC in shallow groundwater can also volatilize and migrate into building structures located above the dissolved phase plume of VOC.

Migration of metals and PAH compounds through site soil and sediment and into on-site buildings is not expected due to the generally low mobility of metals and PAH compounds in these media.

6.5 <u>Evaluation for Immediate Response Actions</u>

Immediate response actions regarding the identified releases at the site were not warranted for the following reasons:

- Site conditions that would warrant a "Two Hour" notification was not identified for the releases.
- A "72 Hour" notification was not identified for the releases.
- A Substantial Release Migration was not identified.
- Accelerated response actions were not necessary to prevent, eliminate, or minimize damage to health, safety, public welfare, or the environment.

7.0 NUMERICAL RANKING SYSTEM (NRS) SITE SCORING —TIER CLASSIFICATION

NRS Site Scoring was conducted in accordance with 310 CMR 40.1500. Details of the site scoring are presented on MADEP forms BWSC107 and BWSC107A, copies of which are included in Appendix E. The NRS score for the disposal site was 212. In accordance with



the MCP, the disposal site is Tier Classified as Tier II. In accordance with Tier Classification requirements, the Chief Municipal Officer and Local Board of Health of the Town of Franklin have been notified of this submittal, and a Legal Notice of a Tier Classification has been made in accordance with 310 CMR 40.1403. Copies of these notifications are included in <u>Appendix F</u>.

8.0 CONCLUSIONS

Based upon the results of this assessment, the following conclusions were drawn:

Soil

- A release of TCE and PCE to surficial soil was identified near the northern boundary of the disposal site. The source of the release was attributed to historic surficial releases of chlorinated solvents.
- A release of TCE and PCE to surficial and vadose zone soil was identified near the intersection of the Lot 27 building and Mine Brook (soil boring B-4). The source of the release in this area may be attributed to historic surficial releases of chlorinated solvents and to a potential subsurface source of VOCs in groundwater.
- A release of beryllium and lead in surficial soil was identified. The presence of metals at levels above applicable soil standards may be attributed to natural sources and/or historic filling practices at the site.

Shallow Overburden Groundwater

- A release of TCE and PCE to shallow groundwater was identified in the area
 of the surficial soil release of chlorinated solvents. The release of TCE and
 PCE in shallow groundwater may be attributed to the migration of
 chlorinated solvents spilled onto surface soil to shallow groundwater.
- A release of lead in shallow groundwater was identified, and may be attributed to leaching of metals to groundwater from fill deposited at the site.

Bedrock Groundwater

 A release of TCE and PCE to bedrock groundwater was identified at monitoring well MW-13 and MW-14. The release may be attributed to a potential additional source from within or beneath the on-site buildings. Additional investigation following the demolition of site buildings may be warranted to evaluate the source of VOCs in deep groundwater.

Sediment

- The presence of PAH in Mine Brook sediment along the western boundary of the subject site may be attributed to historic filling associated with the channelizing of Mine Brook or to historic discharges from the former manufacturing facility.
- No releases to surface water or to environmental media surrounding the former UST were identified during site investigation activities conducted by Fuss & O'Neill.



- Comprehensive Response Actions are necessary at the disposal site, including Tier Classification pursuant to 310 CMR 40.0500. Based on the information documented in this report and in the attached transmittals, the disposal site qualifies as a Tier II Site. The following transmittal documents are included in <u>Appendix E</u> of this report:
 - o BWSC107 Tier Classification Transmittal Form
 - BWSC107A NRS Scoresheet
 - BWSC107B Tier II Compliance History

9.0 CONCEPTUAL PHASE II SCOPE OF WORK

A conceptual Phase II Scope of Work has been prepared in support of the submittal of the MCP Phase I and Tier II Classification document presented herein, pursuant to the MCP (310 CMR 40.0510).

9.1 Assessment Plan

The anticipated assessment plan includes the additional soil and groundwater assessment subsequent to the demolition of the Lot 27 building. The assessment may include additional soil borings and monitoring wells in the vicinity of the identified releases as well as within the footprint of the building.

Based on the apparent limited extent of PAH compounds in sediment, it is anticipated that additional sampling may be conducted concurrent with the performance of remedial response actions in the area of sediment sample SD-01.

9.2 <u>Projected Schedule</u>

As detailed herein, a significant proportion of Phase II assessment activities have been conducted. The additional assessment activities will be conducted following demolition of the Lot 27 building. A schedule for demolition has not been currently established by the Town. It is understood that assessment response actions shall be implemented in accordance with the timelines specified in the MCP.

9.3 <u>Projected Cost</u>

The projected cost for comprehensive site assessment activities outlined in <u>Section 9.1</u> above is estimated to be in the range of \$25,000 to \$50,000. If further subsurface investigations or alternative risk assessment (Method 3 or Ecological Risk Assessment) are warranted to characterize the site, additional project costs will be necessary. It will not be feasible to conduct the outlined assessment tasks without removal of the existing Lot 27 building. The cost for demolition and associated remediation response actions is estimated to be \$200,000 to \$300,000. The Town has not identified a source for this funding. The Town has applied for funding support from USEPA, and will continue to evaluate alternative funding mechanisms.



The scope of work, budget, and projected schedule are presented based on conceptual agreement and authorization by the potentially responsible party.



10.0 REFERENCES

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TABLES

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT



Table 6 Summary of Soil Analytical Data

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

Sample Location		B 02	B 02	B 04	B 04	R 05	R 05	R 06	B 06	R 10	R 10	R 11	P 11	R 12	R 15	M/M/ 01	M/M/ 01	MW 02	M/M/ 0.2	MW 03	VVVV US	MW OA	M/M/ O4	MW OA	M/M/ OS	MW 05	M/M/ 12	M/M/ 17	M/M/ 17	M/M/ 17
Sample ID		841061130-03	841061130-04	841061130-07	841061130-08	841061130-09	841061130-10	841061130-11	841061130-12	841061201-21	841061201-22	841061201-23	841061201-24	9/10/1201 25	9/1071101 03	841061130-01	841061130-02	841061130-05	841061130-06	841061130-13	841061130-14	841061130-15	841061130-16	841061130-17	841061201-19	841061201-20	841071031-01	841071101-04	841071101-05	841071101-06
Date Collected		11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	12/1/2006	12/1/2006	12/1/2006	12/1/2006	12/1/2006	11/1/2007	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	12/1/2006	12/1/2006	10/31/2007	11/1/2007	11/1/2007	11/1/2007
Sample Type	UNITS	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Duplicate 1	Primary	Primary	Primary	Primary	Primary	Primary	Duplicate 1						
	feet	r minary	r minary	rilliary	Filliary		r illiary	rilliary 0	1 1	r i i i i i i i i	r i i i i i i i	r i i i i i i i	r i i i i i i i	r i i i i i i i	riiilai y		2	rilliary	Filliary	rilliary	r i i i i i i i	r i i i i i i i i	Duplicate i	r i i i i i i i		r minary	10		7 TITTIAL Y	2 Duplicate 1
Starting Depth	feet	0	3	0	3	0.4	3	0	1	0	3	0	7	0	Z	0.5	3	0	3	0	3	0	0	3	0.4	3	10	0.3	0	0
Ending Depth	reet	2	/	2	0	2	/	0.5	2	2	1	2	/	Z	4	2	3	3	/	2	1	2	2	/	2	/	12	2	8	8
Metals (EPA Method 6010)																														
Antimony	mg/kg	ND < 0.56	ND < 0.59	ND < 0.57	ND < 0.57	ND < 0.53	ND < 0.54	ND < 0.54	ND < 0.59	ND < 0.50	ND < 0.50	ND < 0.54	ND < 0.60	ND < 0.56	NA	ND < 0.53	ND < 0.56	ND < 0.59	ND < 0.57	ND < 0.55	ND < 0.55	ND < 0.56	ND < 0.53	ND < 0.55	6.5	6.9	NA	1.0	ND < 0.18	ND < 0.18
Arsenic	mg/kg	1.2	ND < 0.59	ND < 0.57	2.0	1.1	ND < 0.54	ND < 0.54	1.8	ND < 0.50	ND < 0.50	1.8	ND < 0.60	3.1	NA	ND < 0.50	ND < 0.56	6.6	2.6	ND < 0.55	ND < 0.55	ND < 0.56	ND < 0.53	ND < 0.55	3.1	ND < 0.50	NA	3.0	1.3	0.84
Barium	mg/kg	36	20	48	24	39	10	28	26	16	9.6	23	34	30	NA	16	24	36	36	18	11	14	9.2	17	110	55	NA	26	15	15
Beryllium	mg/kg	0.21	0.19	0.34	0.36	0.25	0.13	0.18	0.16	0.7	0.91	0.16	0.38	0.26	NA	0.19	0.57	0.22	0.15	0.12	0.17	0.24	0.081	0.15	0.37	0.17	NA	0.35	0.24	0.21
Cadmium	mg/kg	0.17	0.14	0.34	0.19	0.22	ND < 0.11	0.46	0.13	ND < 0.10	ND < 0.10	0.26	ND < 0.12	0.19	NA	ND < 0.10	ND < 0.11	0.13	ND < 0.11	0.16	ND < 0.11	0.15	ND < 0.11	ND < 0.11	0.54	0.18	NA	0.6	0.37	0.17
Chromium	mg/kg	7.1	6.0	8.4	5.4	5.1	3.5	5.8	7.4	5.2	1.9	5.4	4.4	6	NA	3.2	5.3	35	4.1	2.2	5.5	6.0	1.4	5.7	27	26	NA	24	15	3.8
Copper	mg/kg	91	43	13	18	32	3.4	31	20	6.3	1.9	8.5	2.9	37	NA	4.9	12	160	9.0	5.0	2.9	2.0	2.5	25	29	9.5	NA	110	11	3.1
Lead	mg/kg	40	18	8.4	22	20	1.6	97	25	2.9	4.8	17	4.3	93	NA	4.9	8.1	25	89	9.2	2.6	3.4	1.5	4.7	780	310	NA	68	7.1	2.5
Mercury (EPA Method 7471)	mg/kg	0.029	ND < 0.024	0.034	0.051	0.023	ND < 0.022	ND < 0.021	0.065	0.023	ND < 0.021	0.032	ND < 0.024	0.044	NA	ND < 0.021	ND < 0.022	0.14	ND < 0.023	ND < 0.022	ND < 0.022	ND < 0.023	ND < 0.021	ND < 0.022	0.073	ND < 0.023	NA	0.12	0.028	ND < 0.024
Nickel	mg/kg	4.0	3.6	23	37	4.9	14	10	2.6	3.6	1.0	3.2	1.7	130	NA	2.6	3.3	6.2	5.0	3.2	1.8	2.0	6.5	2.0	6.4	6.3	NA	4.3	4.7	2.2
Zinc	mg/kg	85	63	20	26	48	6.8	71	14	22	15	48	8.4	28	NA	10	13	27	54	14	6.3	4.0	4.2	16	310	84	NA	73	28	11
VOC (EPA Method 8260)																														
1,1,1-trichloroethane	μg/kg	ND < 5.1	ND < 5.8	ND < 570	ND < 1,100	ND < 5.0	ND < 5.3	ND < 1,100	ND < 5.8	ND < 270	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	73	17	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Acetone	μg/kg	ND < 20	ND < 23	ND < 2,300	ND < 4,500	ND < 20	ND < 21	ND < 4,300	ND < 23	ND < 1,100	ND < 22	ND < 20	30	ND < 21	ND < 1,200	ND < 21	ND < 22	ND < 21	ND < 22	ND < 21	ND < 22	ND < 21	ND < 20	ND < 18	ND < 18	ND < 22	ND < 3.2	11	28	35
M/P-xylenes	μg/kg	7.0	ND < 5.8	ND < 350	ND < 340	ND < 5.0	ND < 5.3	ND < 300	ND < 5.8	ND < 270	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	ND < 5.2	ND < 4.9	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Methyl ethyl Ketone	μg/kg	ND < 10	ND < 12	ND < 1,100	ND < 2,300	ND < 10	ND < 10	ND < 2,100	ND < 12	ND < 550	ND < 11	ND < 10	ND < 11	ND < 10	ND < 1,200	ND < 10	ND < 11	ND < 11	ND < 11	ND < 10	ND < 11	ND < 10	ND < 9.8	ND < 9.2	ND < 9.2	ND < 11	ND < 3.2	ND < 6.4	6.6	7.4
Naphthalene	μg/kg	ND < 5.1	10	ND < 110	ND < 110	ND < 5.0	ND < 5.3	ND < 110	ND < 5.8	ND < 100	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	260	ND < 5.4	ND < 5.2	2300	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	140	ND < 4.7	ND < 5.2
Tetrachloroethylene	μg/kg	ND < 5.1	ND < 5.8	15,000	20,000	110	22	34,000	310	4,300	48	40	45	11	40,000	ND < 5.3	ND < 5.6	28	45	130	120	13	18	26	ND < 4.6	ND < 5.6	6.4	ND < 6.4	ND < 4.7	ND < 5.2
Toluene	μg/kg	17	ND < 5.8	ND < 350	ND < 340	ND < 5.0	ND < 5.3	ND < 300	ND < 5.8	ND < 270	ND < 5.4	16	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	ND < 5.2	ND < 4.9	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Trichloroethylene	μg/kg	ND < 5.1	ND < 5.8	19,000	31,000	58	9.6	6,700	79	9,300	150	5	ND < 5.7	6.5	9,200	ND < 5.3	ND < 5.6	12	21	150	67	37	44	24	ND < 4.6	ND < 5.6	3.5	ND < 6.4	ND < 4.7	ND < 5.2
EPH with Targets (MADEP Method)																														
C11-C22 Aromatics	μg/kg	17,000	32,000	40,000	61,000	33,000	ND < 11,000	92,000	ND < 12,000	25,000	ND < 10,000	18,000	ND < 11,000	28,000	NA	100,000	16,000	110,000	20,000	52,000	81,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	ND < 11,000	NA	60,000	17,000	ND < 12,000
C19-C36 Aliphatics	μg/kg	ND < 11,000	ND < 11,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	24,000	ND < 12,000	ND < 10,000	ND < 10,000	ND < 11,000	ND < 11,000	16,000	NA	40,000	14,000	ND < 11,000	ND < 11,000	ND < 10,000	38,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	ND < 11,000	NA	18,000	ND < 12,000	ND < 12,000
2-Methylnaphthalene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	210	ND < 110	200	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	ND < 130	ND < 120	ND < 120
Acenaphthene	μg/kg	740	ND < 110	ND < 110	ND < 110	220	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	ND < 110	ND < 110	560	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	150	ND < 120	ND < 120
Acenaphthylene	μg/kg	240	340	230	250	ND < 100	ND < 110	260	ND < 120	200	ND < 100	ND < 110	ND < 110	120	NA	270	ND < 100	1,300	350	150	120	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	260	ND < 120	ND < 120
Anthracene	μg/kg	ND < 110	320	200	ND < 110	970	ND < 110	340	ND < 120	250	ND < 100	ND < 110	ND < 110	ND < 110	NA	390	ND < 100	2,000	ND < 110	1,800	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	640	ND < 120	ND < 120
Benzo(a)anthracene	μg/kg	ND < 110	1,000	330	310	1,100	ND < 110	560	ND < 120	1,500	ND < 100	ND < 110	ND < 110	490	NA	1,200	ND < 100	4,400	200	2,800	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,800	ND < 120	ND < 120
Benzo(a)pyrene	μg/kg	ND < 110	1,000	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	2,000	ND < 100	ND < 110	ND < 110	ND < 110	NA	1,100	ND < 100	3,900	ND < 110	2,200	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,600	ND < 120	ND < 120
Benzo(b)fluoranthene	μg/kg	ND < 110	1,400	560	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	2,000	ND < 100	ND < 110	ND < 110	290	NA	1,700	ND < 100	5,600	ND < 110	2,600	230	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,800	ND < 120	ND < 120
Benzo(k)fluoranthene	μg/kg	ND < 110	540	170	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	860	ND < 100	ND < 110	ND < 110	ND < 110	NA	600	ND < 100	1,900	ND < 110	1,100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,300	ND < 120	ND < 120
Chrysene	μg/kg	ND < 110	1,200	640	590	1,200	ND < 110	810	ND < 120	120	ND < 100	ND < 110	ND < 110	ND < 110	NA	1,600	ND < 100	4,500	280	2,800	440	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,700	ND < 120	ND < 120
Fluoranthene	μg/kg	490	2,100	380	410	2,300	ND < 110	940	ND < 120	1,300	ND < 100	ND < 110	ND < 110	1,300	NA	2,200	ND < 100	8,200	320	6,700	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	3,900	180	ND < 120
Fluorene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	230	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	390	ND < 110	630	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	280	ND < 120	ND < 120
Indeno (1,2,3-cd)pyrene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	260	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,200	ND < 120	ND < 120
Phenanthrene	μg/kg	350	1,100	150	190	2,300	ND < 110	470	ND < 120	260	ND < 100	ND < 110	ND < 110	940	NA	990	ND < 100	3,500	200	7,600	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	2,800	290	ND < 120
Pyrene	μg/kg	ND < 110	2,000	470	410	2,000	ND < 110	990	ND < 120	1,600	ND < 100	ND < 110	ND < 110	1,200	NA	2,200	ND < 100	6,200	310	6,000	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	3,600	180	ND < 120
VPH with Targets (MADEP Method)																														
C9-C12 Aliphatics	μq/kq	ND < 6,900	ND < 8,800	ND < 7,000	ND < 6,900	ND < 5,900	ND < 6,500	ND < 5,900	ND < 7,800	ND < 6,400	ND < 6,100	ND < 12,000	ND < 8,000	ND < 6,800	NA	ND < 6,200	ND < 6,500	ND < 7,900	ND < 7,000	ND < 6,400	ND < 6,400	ND < 7,600	ND < 6,000	ND < 6,100	14,000	ND < 7,600	NA	ND < 6,400	ND < 6,100	ND < 6,000

NOTES:
--- Not applicable
ND < X: Compound not detected above laboratory reporting limit
NA: Not analyzed
NE: Not established
VOC: Volatile organic compounds
EPH: Extractable Petroleum Hydrocarbons
Bold and color-shaded values indicate exceedence of one or more regulatory criteria listed on Table 11.



Table 6 (Continued) MADEP Criteria for Detected Compounds in Soil and Sediment

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

				Regulatory Ci	riteria for Soil			MADEDCH
		MA Method 1 S-1	MA Method 1 S-1	MA Method 1 S-2	MA Method 1 S-2	MA Method 1 S-3	MA Method 1 S-3	MADEP Stage I Freshwater Sediment
		Std Application	Std Application	Std Application	Std Application	Std Application	Std Application	Screening Criteria*
		for GW-2 area	for GW-3 area	for GW-2 area	for GW-3 area	for GW-2 area	for GW-3 area	Screening Criteria
Total Metals (via Method 6010/7471)	UNITS							
Antimony	mg/kg	20	20	30	30	30	30	NE
Arsenic	mg/kg	20	20	20	20	20	20	33
Barium	mg/kg	1,000	1,000	3,000	3,000	5,000	5,000	NE
Beryllium	mg/kg	1	1	1	1	3	3	NE
Cadmium	mg/kg	2	2	30	30	30	30	5.0
Chromium	mg/kg	30	30	200	200	200	200	110
Copper	mg/kg							150
Lead	mg/kg	300	300	300	300	300	300	130
Mercury	mg/kg	20	20	30	30	30	30	0.18
Nickel	mg/kg	20	20	700	700	700	700	49
Thallium	mg/kg	8,000	8,000	60,000	60,000	80,000	80,000	NE
Zinc	mg/kg	2,500	2,500	3,000	3,000	5,000	5,000	460
VPH (MADEP Method)	,	4 000 000	4 000 000	0.500.000	0.500.000	5 000 000	5 000 000	
C9-C12 Aliphatics	μg/kg	1,000,000	1,000,000	2,500,000	2,500,000	5,000,000	5,000,000	
VOC (Method 8260B)								
1,1,1-trichloroethane	μg/kg	500,000	500,000	600,000	1,000,000	600,000	3,000,000	NE
Acetone	μg/kg	60,000	60,000	60,000	60,000	60,000	60,000	NE
M/P-xylenes	μg/kg	300,000	300,000	5,000,000	5,000,000	300,000	300,000	NE
Methyl ethyl Ketone	μg/kg	40,000	40,000	40,000	40,000	40,000	40,000	NE
Naphthalene	μg/kg	40,000	500,000	2,000,000	2,000,000	40,000	3,000,000	NE
Tetrachloroethylene	μg/kg	10,000	30,000	10,000	200,000	10,000	1,000,000	NE
Toluene	μg/kg	300,000	500,000 90.000	2,500,000 2,000	2,500,000 700.000	300,000 2.000	1,000,000 2,000,000	NE NF
Trichloroethylene	μg/kg	2,000	90,000	2,000	700,000	2,000	2,000,000	INE
EPH (MADEP Method)	ua /lea	2.500.000	2.500.000	5.000.000	5.000.000	5.000.000	5.000.000	NE
C19-C36 Aliphatics C11-C22 Aromatics	μg/kg	800.000	800,000	2.000,000	2.000,000	5,000,000	5,000,000	NE NF
	μg/kg	500,000	500,000	1.000.000	1.000.000	2.000,000	1.000.000	NE NF
2-Methylnaphthalene Acenaphthene	μg/kg μg/kg	1.000.000	1.000.000	2,500,000	2,500,000	5,000,000	4.000.000	NE NE
Acenaphthylene	μg/kg μg/kg	1,000,000	100.000	2,500,000	1.000.000	2,500,000	1,000,000	NF NF
Aceraphthylene	μg/kg μg/kg	1.000.000	1.000.000	2,500,000	2,500,000	5,000,000	5.000.000	57
Benzo(a)anthracene	μg/kg	7.000	7,000	40.000	40.000	300.000	300.000	110
Benzo(a)pyrene	μg/kg	2,000	2,000	4,000	4.000	30.000	30.000	150
Benzo(b)fluoranthene	μg/kg	7.000	7.000	40.000	40.000	300.000	300.000	NF
Benzo(ghi)perylene	μg/kg	1.000.000	1.000.000	2,500.000	2,500,000	2,500,000	2.500,000	NF
Benzo(k)fluoranthene	μg/kg	70.000	70.000	400.000	400.000	3,000,000	3,000,000	NE NE
Chrysene	μg/kg	7,000	7,000	10.000	10.000	40.000	40.000	170
Dibenzo(a,h)anthracene	μg/kg μg/kg	700	7,000	4.000	4.000	30.000	30.000	33
Fluoranthene	μα/kg	1.000.000	1.000.000	3.000.000	3.000.000	5.000.000	5,000,000	420
Fluorene	μg/kg	1,000,000	1,000,000	3,000,000	2.000,000	5,000,000	4.000.000	77
Indeno (1,2,3-cd)pyrene	μg/kg	7.000	7.000	40.000	40.000	300.000	300.000	NE NE
Naphthalene	μg/kg	40,000	500,000	2,000,000	2,000,000	40,000	3,000,000	180
Phenanthrene	μg/kg	1,000,000	100,000	2,500,000	100,000	2,500,000	100,000	200
Pyrene	μg/kg	1,000,000	1,000,000	3,000,000	3,000,000	5,000,000	5,000,000	200

NOTES:

μg/kg: micrograms per kilogram mg/kg: milligrams per kilogram NE: not established

S: soil * Sedim GW: groundwater of Ma VPH: Volatile Petroleum Hydrocarbons by Ma

VOC: volatile organic compounds
EPH: Extractable Petroleum Hydrocarbons
MADEP: Massachusetts Department of Environmental Protection
**Sediment Screening Critical Incorporate Threshold Effect Concentral

*Sediment Screening Criteria incorporate Threshold Effect Concentrations of MacDonald et al. (2000) and revised Sediment Screening Criteria published

by MADEP (2005).



Table 9 Summary of Groundwater Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

Sample Location		MCP Regulat	ory Standards	MW-01	MW-01	MW-01	MW-02	MW-02	MW-03	MW-03	MW-04	MW-05	MW-05	MW-13	MW-13	MW-14	MW-16	MW-17	MW-17
Sample ID		MA Method 1 GW	MA Method 1 GW	841061208-27	841061208-28	841071106-03	841061208-30	841071107-10	841061208-32	841071107-11	841061208-29	841061208-31	841071107-09	841071106-05	841071106-06	841071106-04	841071107-08	841071106-01	841071106-02
Date Collected		Std Application	Std Application	12/8/2006	12/8/2006	11/6/2007	12/8/2006	11/7/2007	12/8/2006	11/7/2007	12/8/2006	12/8/2006	11/7/2007	11/6/2007	11/6/2007	11/6/2007	11/7/2007	11/6/2007	11/6/2007
Sample Type		for GW-2 area	for GW-3 area	Primary	Duplicate	Primary	Duplicate	Primary	Primary	Primary	Duplicate								
Groundwater Parameters	UNITS																		
pH	SU			5.97	5.97	6.35	6.59	6.86	6.33	5.91	5.97	NA	6.03	5.68	5.68	5.78	5.91	6.36	6.36
Specific Conductance	μS/cm			464	464	470	1,727	1,547	1,534	3,129	2,010	NA	NA	3,761	3,761	3,065	2,543	494	494
Temperature	C deg			13.1	13.1	17	8.3	13.7	7.6	14.1	10.5	NA	13.2	12	12	15.5	14.9	17.9	17.9
Turbidity	ntu			36	36	20.5	50	NA	500	NA	15	NA	NA	4.64	4.64	3.3	8.9	4.4	4.4
Dissolved Oxygen	mg/l			0.4	0.4	0.1	7.7	4.9	6.9	4.9	2.6	NA	NA	2.8	2.8	3.5	1.6	0.1	0.1
ORP	mv			-35.0	NA	-116.2	59	45.8	93.2	51.3	45.1	NA	NA	30	30	20.2	52.6	-53.2	NA
Metals (EPA Method 6010)																			
Barium, Total	mg/l	NE	50	0.042	0.038	0.031	0.15	NA	0.21	NA	0.14	0.83	NA	0.24	0.25	0.12	0.11	0.06	0.061
Barium, Dissolved	mg/l	NE	50	NA	NA	NA	NA	0.18	NA	0.17	NA	NA	0.39	NA	NA	NA	NA	NA	NA
Beryllium	mg/l	NE	0.05	ND < 0.0010	ND < 0.0010	NA	ND < 0.0010	NA	0.0087	NA	ND < 0.0010	0.0018	NA						
Cadmium	mg/l	NE	0.004	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	NA	ND < 0.0020	NA	ND < 0.0020	0.0034	NA	ND < 0.0020					
Chromium	mg/l	NE	0.3	ND < 0.010	ND < 0.010	0.0029	ND < 0.010	NA	0.036	NA	ND < 0.01	0.092	NA	ND < 0.0020					
Copper	mg/l	NE	NE	ND < 0.010	ND < 0.010	NA	0.015	NA	0.018	NA	ND < 0.01	0.073	NA						
Lead, Total	mg/l	NE	0.01	0.014	0.012	0.0066	ND < 0.0040	NA	0.098	NA	ND < 0.0040	1.9	NA	0.0033	0.0053	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020
Lead, Dissolved	mg/l	NE	0.01	NA	NA	NA	NA	0.0026	NA	0.006	NA	NA	0.094	NA	NA	NA	NA	NA	NA
Nickel	mg/l	NE	0.2	ND < 0.010	ND < 0.010	NA	0.15	NA	0.054	NA	0.017	0.12	NA						
Zinc	mg/l	NE	0.9	0.023	0.015	NA	0.057	NA	0.17	NA	0.028	0.73	NA						
VOC (EPA Method 8260)																			
1,1,1-trichloroethane	μg/l	4,000	20,000	ND < 1.0	1.8	ND < 1.0													
cis-1,2-dichloroethylene	μg/l	100	50,000	ND < 1.0	ND < 1.0	ND < 1.0	ND < 1.0	1.7	ND < 1.0	ND < 1.0	8	ND < 1.0							
Methyl tert butyl ether	μg/l	50,000	50,000	ND < 1.0	1.8	ND < 1.0	ND < 1.0	1.5	1.4	ND < 1.0	ND < 1.0	ND < 1.0	ND < 1.0						
Tetrachloroethylene	μg/l	50	30,000	ND < 1.0	ND < 1.0	ND < 1.0	6.6	23	43	74	240	ND < 1.0	1.3	290	260	12	41	ND < 1.0	ND < 1.0
Trichloroethylene	μg/l	30	5,000	ND < 1.0	ND < 1.0	ND < 1.0	6.6	25	40	59	150	ND < 1.0	ND < 1.0	60	56	20	45	ND < 1.0	ND < 1.0

NOTES:

--- Not applicable
ND <X: Compound not detected above laboratory reporting limit

NA: Not analyzed NE: Not established

µS/cm: microsiemens per centimeter

C deg: degrees Celcius ntu: nephelometric turbidity units ORP: Oxidation-reduction potential

mv: millivolts

mg/l: milligrams per liter µg/l: micrograms per liter

VOC: Volatile organic compounds

Bold and color-shaded values indicate exceedence of one or more regulatory criteria.



Table 11 Summary of Sediment Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

Sample Location		SD-01	SD-02	SD-03	SD-03	SD-04	SD-05*	SD-05*	SD-06*	SD-07*
Sample ID		841070426-06	841070426-07	841070426-08	841070426-09	841070426-10	937071025-01	937071025-02	937071025-04	937071025-03
Date Collected		4/26/2007	4/26/2007	4/26/2007	4/26/2007	4/26/2007	10/25/2007	10/25/2007	10/25/2007	10/25/2007
Sample Type	UNITS	Primary	Primary	Primary	Duplicate 1	Primary	Primary	Duplicate 1	Primary	Primary
Starting Depth	feet	0	0	0	0	0	0	0	0	0
Ending Depth	feet	0.2	0.2	0.2	0.2	0.2	0.25	0.25	0.25	0.25
Metals (EPA Method 6010)										
Arsenic	mg/kg	ND < 0.31	0.75	ND < 0.30	1.2	ND < 0.30	NA	NA	NA	NA
Barium	mg/kg	15	20	22	16	9.0	NA	NA	NA	NA
Beryllium	mg/kg	0.13	0.16	0.16	0.15	0.15	NA	NA	NA	NA
Cadmium	mg/kg	0.14	0.13	0.16	0.14	0.14	NA	NA	NA	NA
Chromium	mg/kg	1.6	1.3	0.75	1.1	2.9	NA	NA	NA	NA
Copper	mg/kg	7.0	6.1	1.8	1.9	3.2	NA	NA	NA	NA
Lead	mg/kg	8.6	5.9	4.8	6.6	13	NA	NA	NA	NA
Nickel	mg/kg	5.4	3.6	0.69	1.5	1.4	NA	NA	NA	NA
Thallium	mg/kg	0.55	0.99	0.69	ND < 0.30	ND < 0.30	NA	NA	NA	NA
Zinc	mg/kg	23	18	15	16	12	NA	NA	NA	NA
VOC (EPA Method 8260)										
Acetone	μg/kg	ND < 5.2	ND < 5.4	ND < 4.6	ND < 5.1	7.8	NA	NA	NA	NA
Tetrachloroethylene	μg/kg	7.6	37	ND < 4.6	ND < 5.1	ND < 4.5	NA	NA	NA	NA
Trichloroethylene	μg/kg	ND < 5.2	12	ND < 4.6	ND < 5.1	ND < 4.5	NA	NA	NA	NA
EPH/SVOC (MADEP Method/EPA Method 8270)										
C11-C22 Aromatics	μg/kg	14,000	ND < 12,000	ND < 12,000	ND < 12,000	ND < 12,000	NA	NA	NA	NA
C19-C36 Aliphatics	μg/kg	20,000	ND < 12,000	ND < 12,000	ND < 12,000	ND < 12,000	NA	NA	NA	NA
Acenaphthylene	μg/kg	140	ND < 120	ND < 120	ND < 120	ND < 120	77	160	76	170
Anthracene	μg/kg	160	ND < 120	ND < 120	ND < 120	ND < 120	340	280	58	98
Benzo(a)anthracene	μg/kg	330	ND < 120	ND < 120	ND < 120	ND < 120	440	ND < 40	400	920
Benzo(a)pyrene	μg/kg	ND < 120	330	770	350	860				
Benzo(b)fluoranthene	μg/kg	120	ND < 120	ND < 120	ND < 120	ND < 120	470	690	520	1,200
Benzo(k)fluoranthene	μg/kg	140	ND < 120	ND < 120	ND < 120	ND < 120	390	1,000	450	990
Chrysene	μg/kg	ND < 120	480	1,100	500	1,200				
Dibenzo(a,h)anthracene	μg/kg	ND < 120	46	100	ND < 40	160				
Fluoranthene	μg/kg	820	ND < 120	ND < 120	ND < 120	ND < 120	890	2,500	810	2,000
Fluorene	μg/kg	ND < 120	ND < 38	81	ND < 40	ND < 46				
Indeno (1,2,3-cd)pyrene	μg/kg	ND < 120	110	240	130	310				
Phenanthrene	μg/kg	230	ND < 120	ND < 120	ND < 120	ND < 120	ND < 38	1,300	ND < 40	ND < 46
Pyrene	μg/kg	450	ND < 120	ND < 120	ND < 120	ND < 120	740	2,100	950	2,000

NOTES:

--- Not applicable $\ensuremath{\mathsf{ND}}\xspace < \mathsf{X}\xspace$ Compound not detected above laboratory reporting limit

NA: Not analyzed

NE: Not established

VOC: Volatile organic compounds

EPH: Extractable Petroleum Hydrocarbons

SVOC: Semivolatile organic compounds

* Samples collected from SD-05 through SD-07 analyzed for SVOC by EPA Method 8270; other samples analyzed for EPH by MADEP Method Bold and color-shaded values indicate exceedance of Sediment Screening Criteria listed in Table 7.



Table 13 Summary of Surface Water Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

				SW-01	SW-02	SW-03	SW-03	SW-04
		MA Method 1 GW	US EPA Chronic	841070426-01	841070426-02	841070426-03	841070426-04	841070426-05
		Std Application	Criteria Continuous	04/26/2007	04/26/2007	04/26/2007	04/26/2007	04/26/2007
		for GW-3 area	Concentrations	Primary	Primary	Primary	Duplicate	Primary
Metals (Method 6010)	·		_					
Barium	mg/l	50	NE	0.086	0.085	0.084	0.083	0.083
Copper	mg/l	NE	0.0090	0.0040	0.0023	ND < 0.0020	0.0041	0.0023
Lead	mg/l	0.01	0.0025	ND < 0.0020	ND < 0.0020	0.0033	ND < 0.0020	ND < 0.0020
Zinc	mg/l	0.9	0.12	0.018	0.017	0.017	0.016	0.015
VOC (Method 8260)		Varies	Varies	ND < varies	ND < varies	ND < varies	ND < varies	ND < varies
VPH (MADEP Method)								
Methyl tert-butyl ether (MTBE)	μg/l	50,000	NE	ND < 1.0	1.1	ND < 1.0	1.1	ND < 1.0
EPH (MADEP Method)	μg/l	Varies	Varies	ND < varies	ND < varies	ND < varies	ND < varies	ND < varies

ND <X: Compound not detected above laboratory reporting limit

US EPA: United States Environmental Protection Agency VPH: Volatile petroleum hydrocarbons

VPH: Volatile petroleum hydrocarbons EPH: Extractable petroleum hydrocarbons VOC: Volatile organic compounds

NE: Not established



Table 15 Summary of Confirmation Soil Sample Analytical Data

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for the Town of Franklin, Massachusetts

May 2008

	Sample Location	East Sidewall	West Sidewall	North Sidewall	East Bottom	West Bottom	South Sidewall
	Sample Number	841070502-01	841070502-02	841070502-03	841070502-04	841070502-05	841070502-06
	Sample Depth (fbg)	3.0	4.0	3.0	8.0	9.0	4.0
Metals (Method 6010)							
Arsenic	mg/kg	1.0	0.69	0.69	0.77	0.65	1.3
Barium	mg/kg	30	13	12	16	10	30
Beryllium	mg/kg	0.057	0.057	ND < 0.052	0.064	0.067	ND < 0.052
Cadmium	mg/kg	0.36	0.29	ND < 0.10	0.39	0.13	0.26
Chromium	mg/kg	3.6	1.7	0.72	3.2	1.0	1.2
Copper	mg/kg	32	13	4.3	60	11	18
Lead	mg/kg	46	3.6	1.2	45	3.7	30
Nickel	mg/kg	190	3.3	5.3	280	8.0	30
Zinc	mg/kg	94	13	7.0	120	13	26
EPH (MADEP Method)							
C9-C18 Aliphatics	μg/kg	19,000	ND < 10,000	ND < 10,000	66,000	ND < 10,000	ND < 10,000
C19-C36 Aliphatics	μg/kg	61,000	ND < 10,000	ND < 10,000	140,000	ND < 10,000	ND < 10,000
C11-C22 Aromatics	μg/kg	54,000	ND < 10,000	ND < 10,000	59,000	ND < 10,000	ND < 10,000
Acenaphthylene	μg/kg	380	ND < 100	ND < 100	160	ND < 100	ND < 100
Anthracene	μg/kg	320	ND < 100	ND < 100	170	ND < 100	ND < 100
Benzo[a]anthracene	μg/kg	640	ND < 100	ND < 100	410	ND < 100	ND < 100
Benzo[a]pyrene	μg/kg	1,100	ND < 100	ND < 100	360	ND < 100	ND < 100
Benzo[b]fluoranthene	μg/kg	520	ND < 100	ND < 100	180	ND < 100	ND < 100
Benzo[k]fluoranthene	μg/kg	ND < 100	ND < 100	ND < 100	270	ND < 100	ND < 100
Fluoranthene	μg/kg	2,000	ND < 100	ND < 100	1100	ND < 100	ND < 100
Fluorene	μg/kg	ND < 100	ND < 100	ND < 100	120	ND < 100	ND < 100
Indeno[1,2,3-cd]pyrene	μg/kg	260	ND < 100	ND < 100	ND < 97	ND < 100	ND < 100
Phenanthrene	μg/kg	860	ND < 100	ND < 100	610	ND < 100	ND < 100
Pyrene	μg/kg	1,300	ND < 100	ND < 100	570	ND < 100	ND < 100
VOCs (Method 8260)							
Acetone	μq/kq	8.8	6.0	7.2	ND < 5.0	9.3	7.9
Tetrachloroethene	μq/kq	46	6.5	ND < 6.0	12	ND < 5.2	ND < 4.8
1.1.1-trichloroethane	μg/kg	ND < 5.6	ND < 4.8	ND < 6.0	12	ND < 5.2	ND < 4.8
Trichloroethene	μg/kg	6.9	ND < 4.8	ND < 6.0	11	ND < 5.2	ND < 4.8

NOTES:
ND < X: Compound not detected above laboratory reporting limit
EPH: Extractable petroleum hydrocarbons
VOCs: Volatile organic compounds
fbg : feet below grade
mg/kg : milligrams per kilogram
µg/kg : micrograms per kilogram

Created by Reviewed by

SAH DJPF



Table 16 Groundwater Elevation Measurements for On-Site Monitoring Wells Gauged February and November 2007

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase I Initial Site Investigation Report Prepared for Town of Franklin, Massachusetts

May 2008

Date	Location	Time	Depth to Water (feet from PVC)	Absolute Elevation of PVC ^a (feet)	Groundwater Elevation (feet)		
	MW-1	0943	4.16	100.35	96.19		
	MW-2	1300	7.96	98.54	90.58		
12/4/2006	MW-3	1412	8.18	99.73	91.55		
	MW-4	1130	7.56	98.23	90.67		
	MW-5	1515	8.49	104.47	95.98		
	MW-1 ^b	0930	4.60	100.35	95.75		
	MW-2	0911	8.16	98.54	90.38		
	MW-3	0920	8.07	99.73	91.66		
	MW-4		U	nable to locate MW-4			
11/7/2007	MW-5	1100	9.20	104.47	95.27		
	MW-13	0915	6.99	99.31	92.32		
	MW-14	0926	10.41	104.40	93.99		
	MW-16	0924	7.15	100.81	93.66		
	MW-17 ^b	0920	4.61	100.37	95.76		

^aelevation data from surveys conducted December 4, 2006 and November 7, 2007

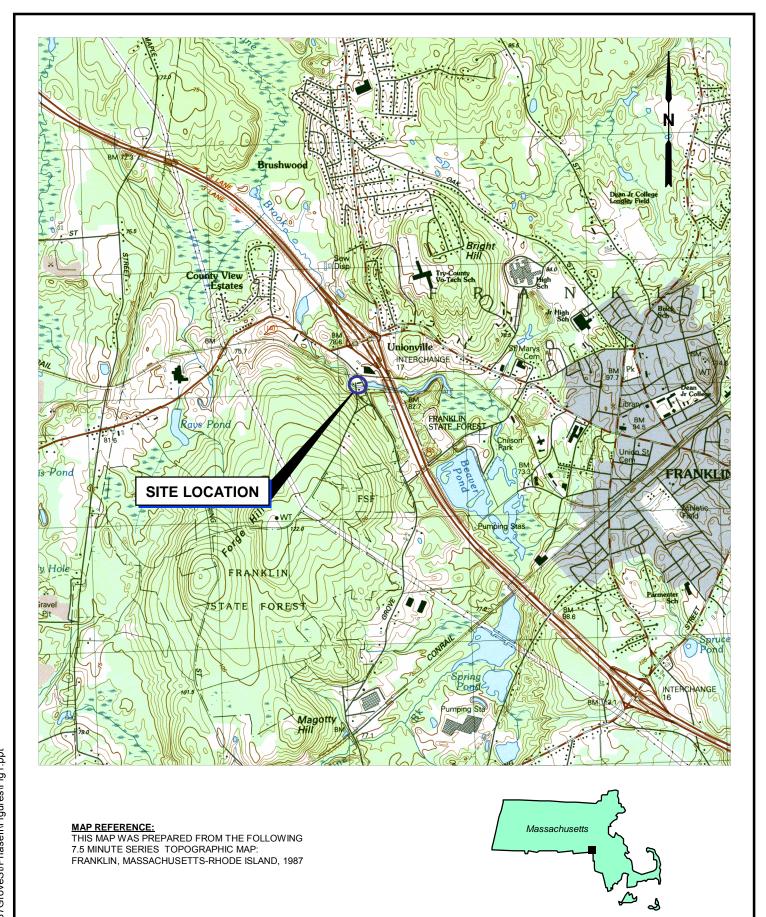
Created by SAH Reviewed by DJPF

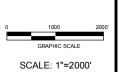
^bMonitoring wells MW-1 and MW-17 were gauged on November 6, 2007 survey data utilize the same arbitrary 100.00-foot benchmark



FIGURES

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT







FOUNDRY CORPORATE OFFICE CENTER
275 PROMENADE ST, SUITE 350, PROVIDENCE RI 02908
401-861-3070 www.FandC

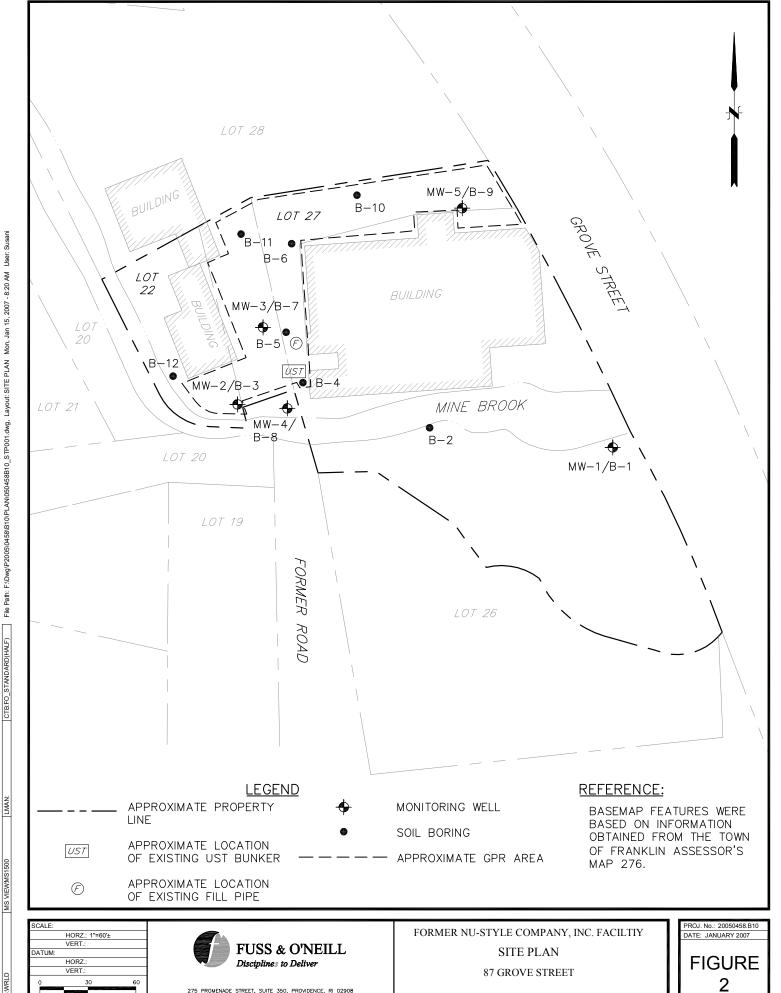
NORFOLK COUNTY, MASSACHUSETTS
SITE LOCATION MAP

FORMER NU-STYLE COMPANY, INC.

87 GROVE STREET FRANKLIN, MASS.

PROJ. No: 20050458.B10 DATE: JANUARY 2007

FIGURE 1



MASSACHUSETTS

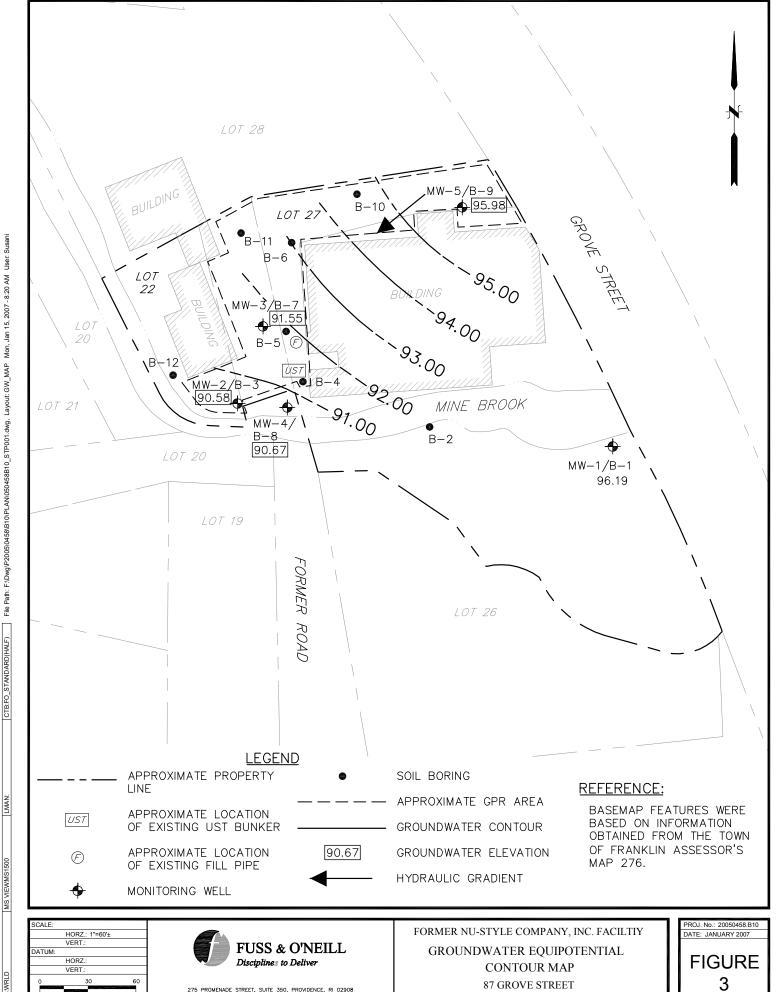
275 PROMENADE STREET, SUITE 350, PROVIDENCE, RI 02908

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FRANKLIN

GRAPHIC SCALE

401.861.3070



MASSACHUSETTS

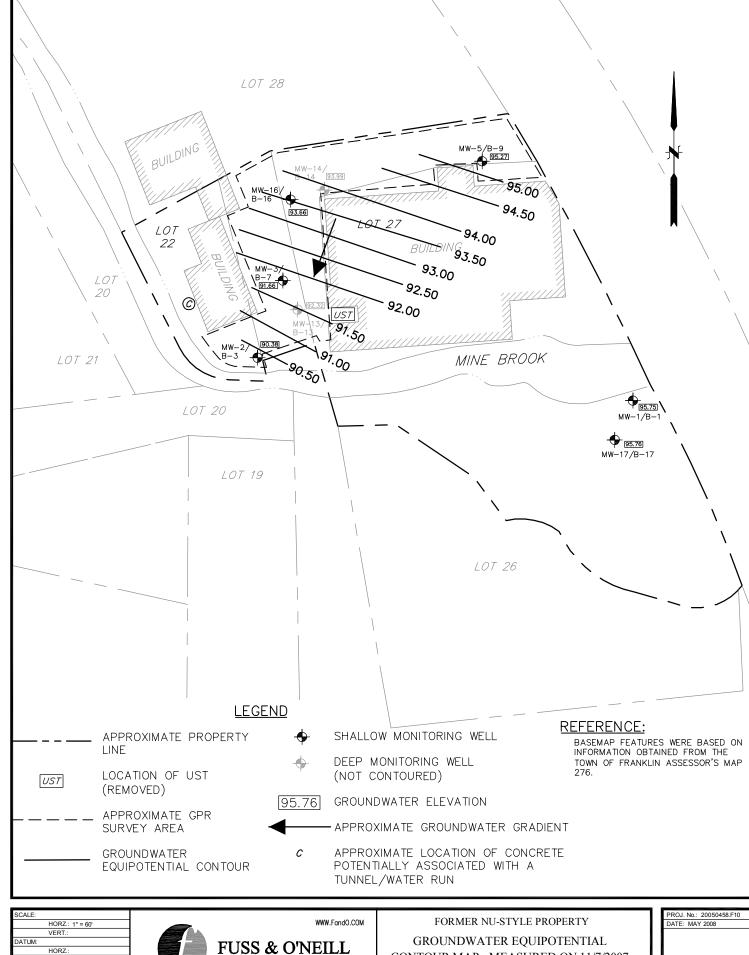
275 PROMENADE STREET, SUITE 350, PROVIDENCE, RI 02908

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FRANKLIN

GRAPHIC SCALE

401.861.3070



HORZ VERT.: GRAPHIC SCALE

MS VIEW

File Path: F: DwgiP2005/0458/B10Nu StyleiPhasel SIR_050458F10_GW_CONTOUR001_NOV2007.dwg, Layout: FIGURE3 - GW CONTOUR MAP Tue, May 13, 2008 - 11:59 AM User: Briank

Disciplines to Deliver

317 IRON HORSE WAY SUITE 204 PROVIDENCE RI 02908 401.861.3070

CONTOUR MAP - MEASURED ON 11/7/2007 87 GROVE STREET

FRANKLIN

MASSACHUSETTS

FIGURE 4



APPENDIX A Limitations of Work Product



LIMITATIONS OF WORK PRODUCT

This document was prepared for the sole use of the Town of Franklin the only intended beneficiaries of our work. Those who may use or rely upon the report and the services (hereafter "work product") performed by Fuss & O'Neill, Inc. and/or its subsidiaries or independent professional associates, subconsultants and subcontractors (collectively the "Consultant") expressly accept the work product upon the following specific conditions.

- 1. Consultant represents that it prepared the work product in accordance with the professional and industry standards prevailing at the time such services were rendered.
- 2. The work product may contain information that is time sensitive. The work product was prepared by Consultant subject to the particular scope limitations, budgetary and time constraints and business objectives of the Client which are detailed therein or in the contract between Consultant and Client. Changes in use, tenants, work practices, storage, Federal, state or local laws, rules or regulations may affect the work product.
- 3. The observations described and upon which the work product was based were made under the conditions stated therein. Any conclusions presented in the work product were based solely upon the services described therein, and not on scientific or engineering tasks or procedures beyond the scope of described services.
- 4. In preparing its work product, Consultant may have relied on certain information provided by state and local officials and information and representations made by other parties referenced therein, and on information contained in the files of state and/or local agencies made available at the time of the project. To the extent that such files which may affect the conclusions of the work product are missing, incomplete, inaccurate or not provided, Consultant is not responsible. Although there may have been some degree of overlap in the information provided by these various sources, Consultant did not attempt to independently verify the accuracy or completeness of all information reviewed or received during the course of this project. Consultant assumes no responsibility or liability to discover or determine any defects in such information which could result in failure to identify contamination or other defect in, at or near the site. Unless specifically stated in the work product, Consultant assumes no responsibility or liability for the accuracy of drawings and reports obtained, received or reviewed.
- 5. If the purpose of this project was to assess the physical characteristics of the subject site with respect to the presence in the environment of hazardous substances, waste or petroleum and chemical products and wastes as defined in the work product, unless otherwise noted, no specific attempt was made to check the compliance of present or past owners or operators of the subject site with Federal, state, or local laws and regulations, environmental or otherwise.
- 6. If water level readings have been made, these observations were made at the times and under the conditions stated in the report. However, it must be noted that fluctuations in water levels may occur due to variations in rainfall, passage of time and other factors



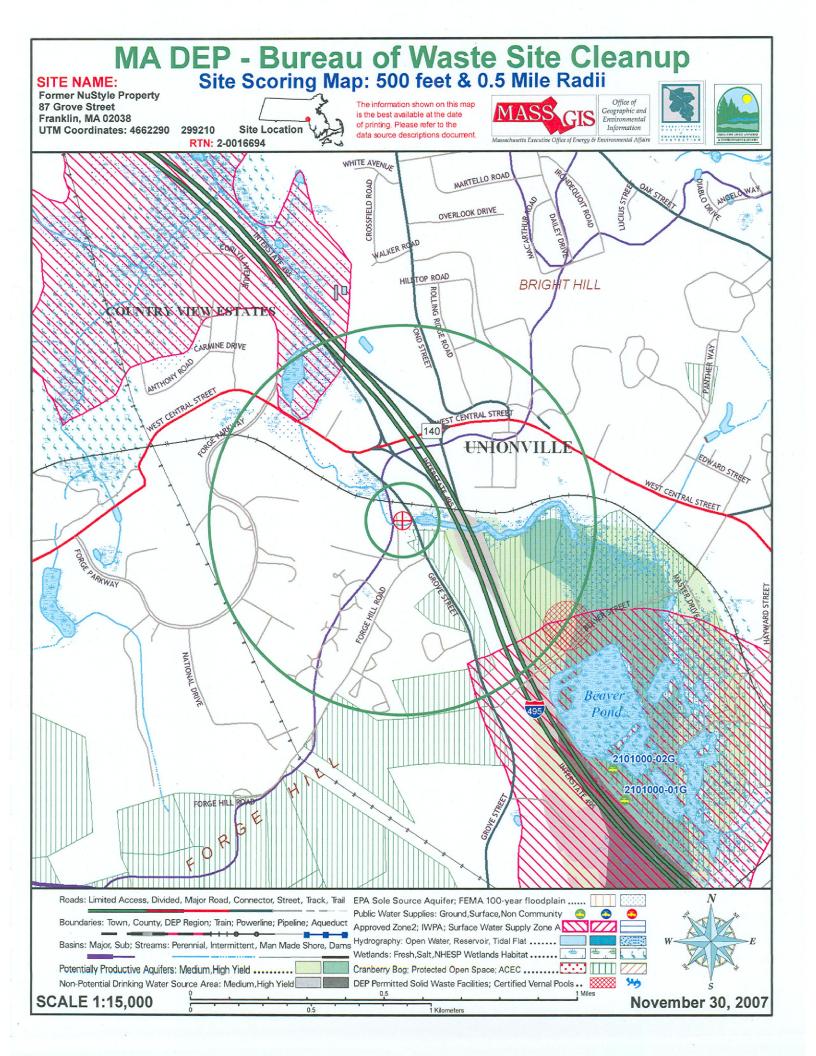
and such fluctuations may effect the conclusions and recommendations presented herein.

- 7. Except as noted in the work product, no quantitative laboratory testing was performed as part of the project. Where such analyses have been conducted by an outside laboratory, Consultant has relied upon the data provided, and unless otherwise described in the work product has not conducted an independent evaluation of the reliability of these tests.
- 8. If the conclusions and recommendations contained in the work product are based, in part, upon various types of chemical data, then the conclusions and recommendations are contingent upon the validity of such data. These data (if obtained) have been reviewed and interpretations made by Consultant. If indicated in the work product, some of these data may be preliminary or screening-level data and should be confirmed with quantitative analyses if more specific information is necessary. Moreover, it should be noted that variations in the types and concentrations of contaminants and variations in their flow paths may occur due to seasonal water table fluctuations, past disposal practices, the passage of time and other factors.
- 9. Chemical analyses may have been performed for specific parameters during the course of this project, as described in the work product. However, it should be noted that additional chemical constituents not included in the analyses conducted for the project may be present in soil, groundwater, surface water, sediments or building materials at the subject site.
- 10. Ownership and property interests of all documents, including reports, electronic media, drawings and specifications, prepared or furnished by Consultant pursuant to this project are subject to the terms and conditions specified in the contract between the Consultant and Client, whether or not the project is completed.
- 11. Unless otherwise specifically noted in the work product or a requirement of the contract between the Consultant and Client, any reuse, modification or disbursement of documents to third parties will be at the sole risk of the third party and without liability or legal exposure to Consultant.
- 12. In the event that any questions arise with respect to the scope or meaning of Consultant's work product, immediately contact Consultant for clarification, explanation or to update the work product. In addition, Consultant has the right to verify, at the party's expense, the accuracy of the information contained in the work product, as deemed necessary by Consultant, based upon the passage of time or other material change in conditions since conducting the work.
- Any use of or reliance on the work product shall constitute acceptance of the terms hereof.



APPENDIX B

MA DEP Bureau of Waste Site Cleanup Site Scoring Map



NRS (21e) SCORING MAP DATA SOURCES

AQUIFERS: USGS-WRD/MassGIS, 1:48,000. Automated by MassGIS from the USGS Water Resources Div. Hydrologic Atlas series manuscripts. The definitions of high and medium yield vary among basins. Source dates 1977-1988.

SOLE SOURCE AQUIFERS: US EPA/MA DEP/MassGIS, various scales. EPA defines them as aquifers that are the 'sole or principal source' of drinking water for a given aquifer service area. Last updated May 1996.

NON-POTENTIAL DRINKING WATER SOURCE AREAS: DEP-BWSC (Bureau of Waste Site Cleanup). Those portions of high and medium yield aquifers, which may not be considered as areas of groundwater conducive to the locations of public water supplies. Please refer to the MCP guidelines for the definitions of these areas. Last updated November 2003.

DEP APPROVED ZONE II's: MA DEP, 1:25,000. As stated in 310 CMR 22.02 'that area of an aquifer which contributes water to a well under the most severe pumping and recharge conditions that can be realistically anticipated.' Digitized from data provided to DEP in approved hydrologic engineering reports. Data are updated continually.

INTERIM WELLHEAD PROTECTION AREAS: DEP-DWS (Division of Water Supply), 1:25,000. These polygons represent an interim Zone II for a groundwater source until an actual one is approved by the DEP Division of Water Supply. The radius of an IWPA varies according to the approved pumping rate. Updated in parallel with the Public Water Supplies data.

PUBLIC WATER SUPPLIES: DEP-DWS, 1:25,000. Community and non-community surface and withdrawal points were field collected using Global Positioning System receivers. The attributes were added from the DEP Division of Water Supply database. Continually updated.

HYDROGRAPHY: USGS/MassGIS. 1:25,000 USGS Digital Line Graph (DLG) data modified by MassGIS. Approximately 40% of the data was provided by USGS and MassGIS created the remainder to USGS specifications. Source dates 1977-1997. Last updated February 2005.

DRAINAGE BASINS: USGS-WRD/MassGIS, 1:24,000. Automated by MassGIS from USGS Water Resources Division manuscripts with approximately 2400 auto-basins as interpreted from 1:24,000 USGS audrangle contour lines. 1987-1993. Last update March 1003.

WETLANDS: UMass Amherst RMP/MassGIS, 1:25,000. Includes nonforested wetlands extracted from the 1999 Land Use datalayer, which was photointerpreted from summer CIR photography. Interpretation was not done in stereo. Also includes, in most areas, forested wetlands from USGS Digital Line Graph (DLG) data.

PROTECTED OPEN SPACE: EOEA (Executive Office of Environmental Affairs) MassGIS, 1:25,000. Includes federal, state, county, municipal, non-profit and protected private conservation and outdoor recreation lands. Ongoing updates.

ACECs: DEM, 1:25,000. Areas of Critical Environmental Concern are areas designated by the Secretary of EOEA as having a number of valuable environmental features coexisting. Projects in ACECs are subject to the highest standards of review and performance. Last updated December 2003.

ROADS: USGS/MassGIS/MHD, 1:100,000. MassGIS extracted roads from the USGS Transportation DLG files. MA Highway Dept. updated roads through July 2004. MassGIS and MA DEP GIS group further edited this layer. Numbered routes are part of the state, U.S. or Interstate highway systems.

POLITICAL BOUNDARIES: MassGIS/USGS, 1:25,000. This datalayer was digitized by MassGIS from mylar USGS quads. Source date is approximately 1985.

DEP PERMITTED SOLID WASTE FACILITIES: DEP-DSW (Division of Solid Waste), 1:25,000. Includes only facilities regulated since 1971. Data includes sanitary landfills, transfer stations and recycling or composting facilities. Facility boundaries were compiled or approximate facility point locations drafted onto USGS quadrangles and automated by the DEP Division of Solid Waste. Last updated November 2003.

NHESP ESTIMATED HABITATS OF RARE WETLANDS WILDLIFE: Polygons show estimated habitats for all processed occurrences of rare wetlands wildlife. Data collected by Natural Heritage & Endangered Species Program and compiled at 1:24,000 or 1:25,000 scale. For use with Wetlands Protection Act Only. Effective 2005-2007.

NHESP CERTIFIED VERNAL POOLS: Points show all vernal pools certified by NHESP/MADFW (Fisheries and Wildlife) as of June 30, 1999. Data compiled at 1:24,000 or 1:25,000 scale. Effective 2005-2007.



APPENDIX C

Phase II ESA, Former Nu-Style Company, Inc. Prepared by Fuss & O'Neill September 2007

Phase II Environmental Site Assessment Report Former Nu-Style Company, Inc. 87 Grove Street (Lots 22 & 27) Franklin, MA

January 2007



275 Promenade Street, Suite 350 Providence, RI 02908



PHASE II ENVIRONMENTAL SITE ASSESSMENT REPORT FORMER NU-STYLE COMPANY, INC. 87 GROVE STREET, FRANKLIN, MA

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PHASE II ENVIRONMENTAL SITE ASSESSMENT REPORT FORMER NU-STYLE COMPANY, INC. 87 GROVE STREET, FRANKLIN, MA

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B Premier Laboratory Certificates of Analysis, Fuss & O'Neill Data Verification Narratives and Certifications, and Data Validation Completeness Worksheets



1.0 INTRODUCTION

1.1 <u>Project Overview and Objectives</u>

The County of Norfolk, Massachusetts (Norfolk County) retained Fuss & O'Neill, Inc. (Fuss & O'Neill) to conduct a Phase II Site Assessment (Phase II) at the former Nu-Style Company, Inc. (the site) located at 87 Grove Street in the Town of Franklin, Massachusetts (the Town). This Phase II was conducted as part of the County Hazardous Materials Brownfield Program, funded under a brownfield assessment grant from the United States Environmental Protection Agency (USEPA).

The objective of the Phase II documented herein was to compile environmental information regarding the site through research, inspections, and field investigations. The Phase II scope of work was developed to determine, to the extent possible with the available resources, the absence or presence and, where applicable, the nature and extent of contaminants in environmental media, to facilitate redevelopment planning at the site, and ultimately to return the property to productive use. The Phase II was conducted in accordance with the Quality Assurance Project Plan (QAPP) Revision 1.0 dated October 2006. This Phase II report has been prepared in accordance with the QAPP Revision 1.0.

The Phase II Report presents the findings of the investigations performed, the conclusions drawn based on those findings, and recommendations with respect to further evaluations or other response actions that may be conducted at the site.

1.2 <u>Assessment Planning and Approvals</u>

Prior to the initiation of Phase II field activities, Fuss & O'Neill prepared a Quality Assurance Project Plan (QAPP) for review and approval by the United States Environmental Protection Agency (USEPA). The QAPP (Revision 0.0) was submitted to USEPA in September 2006. In October 2006, Fuss & O'Neill submitted responses to USEPA comments and questions regarding the QAPP in the form of a document titled QAPP Revision 1.0. The final QAPP (Revision 1.0) was formally approved by USEPA on November 6, 2006. The QAPP, developed in accordance with the USEPA Brownfields Quality Assurance Project Plan Guidance Document, detailed the field and analytical scope and quality control procedures to be implemented during the Phase II.

2.0 BACKGROUND

2.1 <u>Site Description</u>

The site was located at 87 Grove Street in Franklin, Massachusetts. The site has been identified as the Town of Franklin Tax Assessor's Map 276, Lot 22 and Lot 27. Lot 22 covers an area of approximately 9,929 square feet. Lot 27 adjoins Lot 22 to the east and is approximately 42,359 square feet in size. The site was acquired via tax title by the Town of Franklin as a result of foreclosure. Details of the site history are presented in the QAPP and Phase I Environmental Site Assessment (Phase I ESA) report, prepared by Fuss & O'Neill in 2006.



A vacant, partially dilapidated two-story building with a footprint of approximately 11,800 square feet is situated on Lot 27, and a vacant one and one-half-story building with a footprint of approximately 4,000 square feet sits on Lot 22. Mine Brook flows westward along the southern side of the Lot 27 building and turns northward to form the western boundary of Lot 22. Mine Brook flows generally northward to the Charles River.

Utilities located on the subject property include a water line located along the right-of-way known as Old Grove Street, and stormwater drainage lines located throughout the property.

Unrestricted access to the subject property was provided via Grove Street and Old Grove Street.

A portion of a United States Geological Survey (USGS) topographic map depicting the location of the site is provided as <u>Figure 1</u>. A site plan, depicting the boundary of the disposal site, is provided as <u>Figure 2</u>.

2.2 <u>Applicable Regulatory Criteria</u>

Under the provisions of 310 CMR 40.0315, a reportable release exists when analyses of soil or groundwater indicate concentrations of oil or hazardous material that exceed applicable Reportable Concentrations. MADEP must be notified of a reportable release within 120 days of the first knowledge that the release meets one or more reportable criteria. Reportable Concentrations applicable to the subject site were determined based on the current and potential future use of the subject site, nearby property use, and groundwater categories applicable to the subject site in accordance with 310 CMR 40.0932.

Reporting category RCS-1 applies to all soil samples obtained from at or within 500 feet of a residential dwelling or residentially-zoned property, school, playground, recreational area or park, or from within the boundaries of a groundwater resource area. Reporting category RCS-2 applies to all soil samples not obtained from category RCS-1 areas.

Residential dwellings were located less than 500 feet east and south of the subject site. Therefore, in accordance with 310 CMR 40.0361, the RCS-1 reporting category was applied to all soil samples obtained on the subject site.

Reporting category RCGW-1 applies to all groundwater samples obtained from within an area defined as a Current Drinking Water Source Area or Potential Drinking Water Source Area. All groundwater samples not obtained from one of these GW-1 areas falls under the RCGW-2 reporting category.

The subject site was not located in a GW-1 area, and therefore, in accordance with 310 CMR 40.0362, the RCGW-2 reporting category was applied to all groundwater samples obtained on the subject site.



2.3 <u>Topography and Geology</u>

The topography of the site was generally flat, except at the banks of the Mine Brook, where the topography dropped steeply to the river bed (USGS, 1987). The regional topography was hilly and generally drained to Mine Brook.

Surficial material at the site was mapped as loamy udorthents, which generally consist of moderately coarse-grained, deep and moderately deep, fairly well-drained soils (USDA, 2006). Fill described as sand, gravel, silt, and, in some cases, wood and brick was observed to depths of up to 12 feet during drilling conducted on the site as part of the Phase II documented herein.

Bedrock beneath the site was mapped as the grayish-pink to greenish-gray, equigranular to slightly porphyritic Dedham Granite (Hermes et al., 1994). Bedrock was not encountered during drilling and was not visible on or in the immediate vicinity of the site.

2.4 Previous Environmental Investigations

2.4.1 IES Site Investigation Activities

Portions of two reports prepared by IES, Inc. (IES) summarizing environmental investigations previously conducted on the subject property and on the parcel adjacent to the south, were reviewed. The results of the investigations documented in these reports are discussed below.

January 1990

In January 1990, IES completed a report of a Chapter 21E Site Evaluation of 87 Grove Street for Home National bank of Milford, Massachusetts. Portions of the report were available for review at the Franklin Health Department.

The IES investigation included the drilling of soil borings and the collection and analyses of soil and groundwater samples on the subject property and on the parcel adjacent to the south of Lot 27 (Lot 26). Note that the map and parcel numbers have changed since the IES investigation, as summarized in the table below.

Pre	vious	Cui	rrent	Comments		
Map	Lot	Map	Lot	Comments		
72	5	276	22	Site		
72	6	276	27	Site		
72	7	276	26	Adjacent south		

IES collected soil and/or groundwater samples from five borings (B-1 through B-5) drilled on the three parcels. A figure provided by IES shows the approximate boring locations; however because the figure is schematic and is not to scale, the precise boring locations could not be determined.



Two of the borings (B-1 and B-2) were drilled adjacent to underground storage tanks located on Lot 26. Borings B-3 and B-5 were situated on the north side of the Lot 27 building, and boring B-4 was advanced in the exterior "barrel area" north of the Lot 22 garage. Field screening indicated the presence of trace concentrations of volatile organic compounds (VOCs) in the soil at borings B-4 and B-5; therefore, soil from the two borings from a depth of approximately five feet below grade was composited into one sample, which was analyzed for VOCs. No VOCs were detected. Groundwater was not encountered at these two boring locations.

Groundwater samples collected from B-1 and B-2 were also analyzed for VOCs, which were not detected. No information regarding sample analysis for soil or groundwater collected from B-3 was reported; therefore, we infer that no samples were analyzed because field screening did not indicate the presence of VOCs.

IES concluded that no releases of hazardous materials or petroleum products had occurred at the subject property; however, it is Fuss & O'Neill's opinion that the IES investigation was not adequate to definitively rule out releases on the subject property.

July 1991

In July 1991, IES collected soil samples from four additional borings (B-1A through B-4A) to assess whether releases associated with underground storage tanks had occurred. As with the 1990 investigation, only portions of the July 1991 report were available for review at the Franklin Health Department. A copy of the report was also available at the Franklin Fire Department, but copies could not be made. A figure depicting the boring locations was not included with the report.

Soils generally consisted of fill containing loam, sand, gravel, and, in some cases, brick and cinders. Fill materials were observed to depths of up to 8.5 feet below grade. Deeper soils consisted of very dense, fine-grained sand, silt, and gravel. Groundwater was encountered at depths of approximately 8.5 to 9 feet. Monitoring wells were installed within the borings to allow for the collection of groundwater samples.

IES identified releases of chlorinated solvents to soil and groundwater at boring location B-4A, which was situated downgradient of USTs at the site and north of Mine Brook. Based on the apparent vertical distribution of VOCs in soil, IES inferred that the presence of VOCs was the result of a surface release.

2.4.2 Phase I ESA Report, May 2006

A Phase I ESA report, prepared by Fuss & O'Neill in May 2006, identified the following recognized environmental conditions (RECs) at the site:

• The site has a long history (at least 90 years) of manufacturing, including textiles and jewelry. Materials used and stored at the site associated with jewelry manufacturing include cyanides, metals, chlorinated solvents, and petroleum products. Additional substances associated with textile manufacturing were also likely used. There is the potential for surface releases to have occurred associated with the use and storage of



these materials. Files indicate that numerous drums of hazardous waste and petroleum products were situated outside of the site buildings.

- At least one underground storage tank appears to be present on the western side of the Lot 27 building. In addition, a heating oil tank reportedly exists in an underground bunker on the same side of the building. As with any underground tank, there is the potential for releases to have occurred associated with leaks or spills.
- A small tunnel containing slow-flowing water is present beneath the Lot 22 building.
 A review of mapping on file at the Town Building Department suggests that the
 tunnel runs, or ran in the past, from Mine Brook and beneath the Lot 27 building to
 the Lot 22 building. There is the potential that the tunnel was used by the former
 woolen mill for direct waste disposal to Mine Brook prior to the realignment of the
 brook in the 1960s.
- Releases of chlorinated solvents to soil and groundwater were identified on Lot 26, which abuts the site to the south. Due to the proximity of this property to the site, there is the potential for releases that occur on this property to adversely affect groundwater quality at the site. Note that this property was owned and occupied by the same entities that owned and operated the facilities at the site; therefore, there is the potential that similar releases have occurred at the site.
- The southern portion of the site contained a pond that was filled circa 1960. The fill appears to have been placed by a municipality. The nature and origin of the fill are not known.

2.4.3 Phase I ESA Report, January 2007

A Phase I ESA report, was prepared by Fuss & O'Neill in accordance with ASTM E1527-05 dated January 2007. The January Phase I ESA report identified the same RECs that were noted in May 2006, summarized above.

3.0 PHASE II ACTIVITIES

Based on the results of the Phase I ESA, a Phase II scope of work was developed to assess the potential presence of Oil or Hazardous Material (OHM) in soil and groundwater at the site, and to determine if a Reportable Condition pursuant to 310 CMR 40.0360 existed at the site. The Phase II scope of work was implemented in accordance with the approved QAPP.



3.1 <u>Field Investigation Activities</u>

3.1.1 Ground Penetrating Radar Survey

In accordance with the QAPP, a ground penetrating radar (GPR) survey was conducted at the site on November 1, 2006. The GPR survey was conducted to evaluate the potential presence of suspected USTs on the subject property.

The following three anomalies indicative of potential USTs were reported in the areas surveyed at the site:

- One anomaly was identified beneath a raised concrete pad located off the southwest corner of the building on Lot 27.
- One anomaly was identified off the northern edge of the Lot 27 building, northeast of the loading dock.
- One anomaly was identified off the northern edge of the Lot 27 building. The anomaly was elongated parallel to the northern edge of the loading dock, and extended past the western end of the loading dock.

The GPR survey was not conducted in portions of the northern end of Lot 26 due to the presence of parked vehicles. In addition, areas immediately adjacent to the buildings that were overgrown with vegetation or filled with debris could not be surveyed.

3.1.2 Soil Sampling

In accordance with the QAPP, a soil sampling program was conducted at the site on November 30, 2006 and December 1, 2006. Twelve soil borings (B-1 through B-12) were advanced throughout the site utilizing direct-push drilling methods. Borings were advanced to a depth of up to 12 feet below grade. Refer to Figure 2 for a map of soil boring locations.

Two soil samples were collected from boring B-1 through B-11; one sample was collected from the 0 to 2 foot depth interval and one sample was collected from vadose zone soil directly above the water table. Due to a lack of adequate soil recovery at depth, one soil sample was collected from the 0 to 2 foot depth interval of boring B-12.

Twenty-four soil samples, including one field duplicate, were submitted to Premier Laboratory, LLC in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant metals plus barium by EPA Methods 6010B and 7471, cyanide by EPA Method 9012, polychlorinated biphenyls (PCBs) by EPA Method 8082, and petroleum hydrocarbons by Massachusetts Department of Environmental Protection (MADEP) Methods Extractable Petroleum Hydrocarbons (EPH) and Volatile Petroleum Hydrocarbons (VPH). One aqueous trip blank was also submitted each day for analysis of VOCs by EPA Method 8260B. Dedicated sampling equipment was employed; therefore, no equipment blank was evaluated.

Soil boring logs, depicting sample recovery amounts, material descriptions, graphic logs, soil codes, and PID soil screening results are attached in <u>Appendix A</u>.



3.1.3 Monitoring Well Installation and Development

In accordance with the QAPP, groundwater monitoring wells were installed in five of the soil borings advanced at the site. The locations of the monitoring wells are depicted on <u>Figure 2</u>. Detailed monitoring well completion reports are included in <u>Appendix A</u>.

In accordance with the QAPP, Fuss & O'Neill surveyed the relative elevations of the newly installed monitoring wells at the site on December 4, 2006. The survey was conducted relative to an assumed arbitrary vertical datum to evaluate the relative elevation and hydraulic gradient of shallow groundwater beneath the site.

In accordance with the QAPP, Fuss & O'Neill developed the newly installed monitoring wells at the site on December 4 and 6, 2006. Development procedures included the repeated purging and surging of groundwater in the wells to remove fine particles and to improve hydraulic communication between the sand filter pack and surrounding soil formation. Low recovery volumes and slow recharge were observed at wells MW-3 and MW-5.

3.1.4 Low Flow Groundwater Sampling

On December 8, 2006, Fuss & O'Neill collected six groundwater samples, including one duplicate sample, from newly installed monitoring wells at the site. Groundwater samples were collected from monitoring wells MW-1, MW-2 and MW-4 utilizing low-flow sampling techniques, in accordance with the QAPP. These techniques included the following:

- A peristaltic pump with dedicated, disposable tubing was utilized to purge groundwater from each well.
- During purging, field parameters including groundwater pH, temperature, and specific conductivity were measured with a multi-meter.
- A groundwater sample was collected when field parameters stabilized to within limits specified in Fuss & O'Neill's Standard Operating Procedures (SOPs).

Low water volume and slow recharge at wells MW-3 and MW-5 prevented the utilization of low-flow field parameter monitoring techniques. Instead, groundwater samples were collected at low flow rates following limited purging of MW-3, and without significant purging at MW-5.

Groundwater samples were submitted to Premier Laboratory for analysis for priority pollutant 13 (PP-13) metals, barium, VOCs, and petroleum hydrocarbons (MADEP EPH and VPH Methods).

One of the groundwater samples collected during this assessment was a duplicate sample collected from monitoring well MW-1, and submitted to the laboratory for analysis of VOCs for quality control purposes. One trip blank was also collected and submitted to Premier Laboratory for analysis of VOCs.



4.0 INVESTIGATION RESULTS

4.1 <u>Soil PID Field Screening Results</u>

During soil boring advancement, soil samples were collected throughout the soil column at each boring location for field screening for the presence of total VOCs with a photoionization detector (PID). Field screening results indicated that total VOCs were not detected in any soil samples. Field screening results are included in soil boring logs attached as <u>Appendix A</u>.

4.2 <u>Soil Analytical Laboratory Results</u>

A summary of soil analytical data is included in <u>Table 1</u>. The complete Premier Laboratory analytical data packages and associated data verification narratives and certifications for each laboratory report are attached in <u>Appendix B</u>.

Laboratory analytical results of soil samples collected from on-site soil borings documented the presence of the following analytes in soil at concentrations above laboratory reporting limits:

Metals (Method 6010)	VOCs (Method 8260)	PAHs (MADEP EPH Method)
Antimony	Naphthalene	2-Methylnaphthalene
Arsenic	Toluene	Acenaphthene
Barium	M/P-xylenes	Acenaphthylene
Beryllium	Acetone	Anthracene
Cadmium	Tetrachloroethene (PCE)	Benzo(a)anthracene
Chromium	Trichloroethene (TCE)	Benzo(a)pyrene
Copper		Benzo(b)fluoranthene
Lead		Benzo(k)fluoranthene
Nickel		Chrysene
Selenium		Fluoranthene
Silver		Fluorene
Thallium		Indeno(1,2,3-cd)pyrene
Zinc		Phenanthrene
Mercury (Method 7471)		Pyrene

4.3 Groundwater Analytical Results

A summary of groundwater analytical data is included in <u>Table 2</u>. The complete Premier Laboratory analytical data packages and associated data verification narratives and certifications for each laboratory report are attached in <u>Appendix B</u>.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of the following analytes in groundwater at concentrations above laboratory reporting limits:

Metals (Method 6010)	VOCs (Method 8260)
Barium	Methyl tert butyl ether (MTBE)



Beryllium	cis-1,2-Dichloroethene
Cadmium	Tetrachloroethene
Chromium	1,1,1-trichloroethane
Copper	Trichloroethene
Lead	
Nickel	
Zinc	

4.4 Surficial and Subsurface Soil Characterization

In general, the soil within soil borings advanced at the site was observed to consist of mainly fine to medium sand, with varying proportions of gravel and silt. Apparent fill material containing wood, brick, coal and/or coal ash was observed in soil borings advanced throughout the site, and was concentrated in the upper five feet of soil.

Two soil horizons consisting predominately of silt were encountered in soil boring MW-05. The upper silty horizon had a minimum thickness of four feet, and occurred in the 5-10 foot depth interval. The lower silty horizon occurred in the 10-12 foot depth interval, had a minimum thickness of 0.8 feet, and may have extended beyond the maximum boring depth.

5.0 EVALUATION OF ANALYTICAL RESULTS

5.1 Data Verification

Procedures and methodologies for the collection and analyses of soil and groundwater samples were performed consistent with the QAPP and the MCP (310 CMR 40.0017). Analytical data were developed and reviewed in accordance with the MADEP's Compendium of Quality Assurance and Quality Control Requirements and Performance Standards for Selected Analytical Methods (the CAM).

Fuss & O'Neill conducted modified Tier I data verification of the field and analytical data resulting from the assessment documented herein. Modified Tier I verification narratives and certifications, signed by the Fuss & O'Neill Quality Assurance/Quality Control Officer, as well as modified Tier I completeness and verification checklists are attached to each Premier Laboratory report in <u>Appendix B</u>.

Presumptive Certainty was obtained for each data set collected as part of the Phase II investigation. Documentation was provided by Premier Laboratory along with narrative summaries (Appendix B).

With the exception of ethylene dibromide and 1,1,2,2-tetrachloroethene, reporting limits were generally low enough to allow for direct comparison to the applicable criteria (RCS-1 and RCGW-2). Limitations of the instrumentation and sample matrices prevent the laboratory from being able to attain a reporting limit of 5 mg/Kg for ethylene dibromide and 1,1,2,2-tetrachloroethene for all samples. Other volatile organic compounds (VOCs) and extractable petroleum hydrocarbon (EPH) target compounds are reported with elevated reporting limits due to dilutions required for quantification of detected compounds. The usability of the data is not anticipated to be affected by these issues.



5.2 <u>Soil</u>

Soil analytical data are summarized in <u>Table 1</u>. In accordance with 310 CMR 40.0361, the RCS-1 reporting category was applied to all soil samples obtained on the subject site.

Tetrachloroethene and trichloroethene were detected at levels in excess of the RCS-1 criteria in soil samples collected from the 0-5 and 5-10 depth intervals of boring B-04, located near the UST area of the site, north of Mine Brook. These same compounds were detected in the 0-5 foot depth-interval samples collected from B-06 and B-10. These results suggest that the maximum vertical extent of tetrachloroethene and trichloroethene in soil at the site may be limited to the upper five feet of soil on the northern end of the site, but may extend to a depth of at least 10 feet below grade in the area downgradient of the UST area, proximal to the northern bank of Mine Brook.

Laboratory analysis of soil samples collected from the 0-5 and 5-10 foot depth intervals of borings B-04, MW-05, and B-10 documented the presence of beryllium, lead and/or nickel at levels in excess of the RCS-1 criteria. Laboratory analysis of a shallow soil sample collected from the 0-5 foot depth interval of boring B-12 documented the presence of nickel, lead, and/or beryllium at levels in excess of the RCS-1 criteria. As previously noted, due to poor soil recovery during boring operations, a deeper soil sample could not be collected from boring B-12.

Laboratory analysis of a soil sample collected from the 0-5 foot depth interval of boring MW-03 documented the presence of the polycylic aromatic hydrocarbons (PAHs) benzo(a)pyrene and fluorene at levels in excess of the RCS-1 criteria. In addition, benzo(a)pyrene was detected at levels in excess of the RCS-1 criteria in a soil sample collected from the 0-5 foot depth interval of boring B-10.

5.3 Groundwater

A summary of groundwater analytical results is included as <u>Table 2</u>. In accordance with 310 CMR 40.0362, the RCGW-2 reporting category was applied to all groundwater samples obtained on the subject site.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of lead (wells MW-1, MW-3 and MW-5), tetrachloroethene (wells MW-3 and MW-4) and trichloroethene (wells MW-3 and MW-4) at levels in excess of the RCGW-2 criteria.

6.0 CONCLUSIONS

6.1 Soil

A comparison of the soil analytical results, documented herein, to the RCS-1 criteria indicated that pursuant to the MCP (310 CMR 40.0361) a reportable condition existed with regard to soil at the subject site.



Laboratory analytical results of soil samples collected from soil borings advanced on-site documented the presence of the following target analytes at concentrations in excess of one or more criteria:

- Metals: beryllium, lead and nickel
- Volatile Organic Compounds: PCE and TCE
- Polynuclear Aromatic Hydrocarbon Compounds: benzo(a)pyrene and fluorene

6.2 Groundwater

A comparison of the groundwater analytical results, documented herein, to the RCGW-2 criteria indicated that pursuant to the MCP, 310 CMR 40.0362, a reportable condition existed with regard to groundwater at the subject site.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of the following target analytes at concentrations in excess of one or more criteria:

Metals: lead

Volatile Organic Compounds: PCE and TCE

7.0 RECOMMENDATIONS

Based upon the results of this assessment, the following response actions are recommended:

- Following the completion of the data validation process, on January 16, 2007 representatives of the Town of Franklin were informed of the presence of levels of metals, PAH compounds, trichloroethene and tetrachloroethene at reportable concentrations. As such, on or before May 16, 2007 appropriate documentation should be submitted to the MADEP reporting the 120-day reportable condition.
- Further assessment of soil and groundwater is recommended to support MCP-related response actions.
- Assessment of surface water and sediment in Mine Brook is recommended to characterize potential impacts from the release at the subject site.



8.0 REFERENCES

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9.0 LIMITATIONS OF WORK PRODUCT

Those who may use or rely upon the report and the services (hereafter "work product") performed by Fuss & O'Neill, Inc. and/or its subsidiaries or independent professional associates, subconsultants and subcontractors (collectively the "Consultant") expressly accept the work product upon the following specific conditions.

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- 3. The observations described and upon which the work product was based were made under the conditions stated therein. Any conclusions presented in the work product were based solely upon the services described therein, and not on scientific or engineering tasks or procedures beyond the scope of described services.
- 4. In preparing its work product, Consultant may have relied on certain information provided by state and local officials and information and representations made by other parties referenced therein, and on information contained in the files of state and/or local agencies made available at the time of the project. To the extent that such files which may affect the conclusions of the work product are missing, incomplete, inaccurate or not provided, Consultant is not responsible. Although there may have been some degree of overlap in the information provided by these various sources, Consultant did not attempt to independently verify the accuracy or completeness of all information reviewed or received during the course of this project. Consultant assumes no responsibility or liability to discover or determine any defects in such information which could result in failure to identify contamination or other defect in, at or near the site. Unless specifically stated in the work product, Consultant assumes no responsibility or liability for the accuracy of drawings and reports obtained, received or reviewed.
- 5. If the purpose of this project was to assess the physical characteristics of the site with respect to the presence in the environment of hazardous substances, waste or petroleum and chemical products and wastes as defined in the work product, unless otherwise noted, no specific attempt was made to check the compliance of present or past owners or operators of the site with Federal, state, or local laws and regulations, environmental or otherwise.
- 6. If water level readings have been made, these observations were made at the times and under the conditions stated in the report. However, it must be noted that fluctuations in water levels may occur due to variations in rainfall, passage of time and other factors



and such fluctuations may effect the conclusions and recommendations presented herein.

- 7. Except as noted in the work product, no quantitative laboratory testing was performed as part of the project. Where such analyses have been conducted by an outside laboratory, Consultant has relied upon the data provided, and has not conducted an independent evaluation of the reliability of these tests.
- 8. If the conclusions and recommendations contained in the work product are based, in part, upon various types of chemical data, then the conclusions and recommendations are contingent upon the validity of such data. These data (if obtained) have been reviewed and interpretations made by Consultant. If indicated in the work product, some of these data may be preliminary or screening-level data and should be confirmed with quantitative analyses if more specific information is necessary. Moreover, it should be noted that variations in the types and concentrations of contaminants and variations in their flow paths may occur due to seasonal water table fluctuations, past disposal practices, the passage of time and other factors.
- 9. Chemical analyses may have been performed for specific parameters during the course of this project, as described in the work product. However, it should be noted that additional chemical constituents not included in the analyses conducted for the project may be present in soil, groundwater, surface water, sediments or building materials at the site.
- 10. Ownership and property interests of all documents, including reports, electronic media, drawings and specifications, prepared or furnished by Consultant pursuant to this project are subject to the terms and conditions specified in the contract between the Consultant and Client, whether or not the project is completed.
- 11. Unless otherwise specifically noted in the work product or a requirement of the contract between the Consultant and Client, any reuse, modification or disbursement of documents to third parties will be at the sole risk of the third party and without liability or legal exposure to Consultant.
- 12. In the event that any questions arise with respect to the scope or meaning of Consultant's work product, immediately contact Consultant for clarification, explanation or to update the work product. In addition, Consultant has the right to verify, at the party's expense, the accuracy of the information contained in the work product, as deemed necessary by Consultant, based upon the passage of time or other material change in conditions since conducting the work.
- 13. Any use of or reliance on the work product shall constitute acceptance of the terms hereof.



TABLES

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT

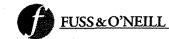


Table 1 Summary of Soil Analytical Data and Objectives

Former Nu-StyleCompany, Inc. Grove Street Franklin, Massachusetts

Phase II Comprehensive Site Assessment Report Prepared for the County of Norfolk, Massachusetts

January 2007

																										,
			MW-01	MW-01	B-02	B-02	MW-02	MW-02	B-04	B-04	B-05	B-05	B-06	B-06	MW-03	MW-03	. MW-04	MW-04	MW-04	MW-05	MW-05	B-10	I B-10 I	1 2044	7.14	T 742
			841061130-01	841061130-02	841061130-03	841061130-04	841061130-05	841061130-06	841061130-07	841061130-08	841061130-09	841061130-10	841061130-11	841061130-12	841061130-13	841061130-14	841061130-15	841061130-16	841061130-17	841061201-19	MW-05 841061201-20	841061201-21	B-10 841061201-22	B-11 841061201-23	B-11 841061201-24	B-12 841061201-25
	•	"MADEP	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	12/01/2006	12/01/2006	12/01/2006	12/01/2006	12/01/2006	12/01/2006	1
		2006	1.25	4.00	1.00	6.00	1.50	6.00	1.00	5.50	1.20	6.00	0.25	. 1.50	1.00	6.00	1.00	1.00	6.00	1.20	6.00	1.00	6.00	1.00	6.00	12/01/2006 1.00
4	1.2	∈RC-S1-	Primary	Primary	Primary	Primary.	Primary.	Primary	Primary.	Primary.	Primary.	Primary	Primary	Primary.	Primary	Primary	Primary	Duplicate 1	Primary	Primary	Primary	Primary	Primary	Primary.	Primary.	Primary
Starting Depth	(feet)	 ′	0.50	3.00	0.00	5.00	0.00	5.00	0.00	5.00	0.40	5.00	0.00	1.00	0.00	5.00	0.00	0.00	5.00	0.40	5.00	0.00	5.00	0.00	5.00	0.00
Ending Depth	(feet)	<u> </u>	2.00	5.00	2.00	7.00	3.00	7.00	2.00	6.00	2:00	7.00	0.50	2.00	2.00	7.00	2.00	2.00	7.00	2.00	7.00	2.00	7.00	2.00	7.00	2.00
Metals (Method 6010)	UNITS	<u> </u>	4	1	<u></u>								,				·	· ·	1	· ·	T	1	((T	, 	
Antimony	mg/kg	20	ND <0.53	ND <0.56	ND <0.56	ND <0.59	ND <0.59	ND <0.57	ND <0.57	ND <0.57	ND <0.53	ND <0.54	ND <0.54	ND <0.59	ND <0.55	ND < 0.55	ND <0.56	ND < 0.53	ND <0.55	6.5	6.9	ND <0.50	ND <0.50	ND <0.54	ND <0.60	ND <0.56
Barium	mg/kg mg/kg	20 1000	ND <0.50	ND <0.56	1.2	ND <0.59	6.6	2.6	ND <0.57	2.0	1.1	ND <0.54	ND <0.54	1.8	ND < 0.55	ND <0.55	ND <0.56	ND <0.53	ND <0.55	3.1	ND <0.50	ND <0.50	ND <0.50	1.8	ND <0.60	3.1
Beryllium	mg/kg mg/kg	0.7	0.19	0.57	0.21	0.19	36	36 0.15	48	24	39	10	28	26	18	11	14	9.2	17	110	55	16	9.6	23	34	30
Cadmium	mg/kg	2	ND <0.10	ND <0.11	0.21	0.19	0.22	0.15 ND <0.11	0.34	0.36	0.25	0.13	0.18	0.16	0.12	0.17	0.24	0.081	0.15	0.37	0.17	0.7	0.91	0.16	0.38	0.26
Chromium	mg/kg	1000	3.2	5.3	7.1	6.0	35	ND <0.11	0.34 8.4	0.19 5.4	0.22 5.1	ND <0.11	0.46 5.8	7.4	0.16	ND <0.11	0.15	ND <0.11	ND <0.11	0.54	0.18	ND <0.10	ND <0.10	0.26	ND <0.12	0.19
Copper	mg/kg	1000	4.9	12	91	43	160	9.0	13	18	32	3.5	5.8 31	20	5.0	5.5	6.0	1.4	5.7	27	26	5.2	1.9	5.4	4.4	6.0
Lead	mg/kg	300	4.9	8.1	40	18	25	89	8.4	22	20	1:6	97	25	9.2	2.6	3.4	2.5	25	29	9.5	6.3	1.9	8.5	2.9	37
Nickel	mg/kg	20	2.6	3.3	4.0	3.6	6.2	5.0	23	37.	4.9	14	10	2.6	3.2	1.8	2.0	6.5	2.0	6.4	6.3	3.6	1.0	3.2	4,3	93
Selenium	mg/kg	400	ND <0.50	NID < 0.56	ND <0.56	ND <0.59	ND <0.59	ND <0.57	·ND <0.57	ND <0.57	ND <0.53	ND <0.54	ND <0.54	ND <0.59	ND <0.55	ND <0.55	ND <0.56	ND <0.53	ND <0.55	ND <0.50	ND <0.50	3.6 ND <0.50	ND <0.50	3.2 ND <0.54	1.7 ND <0.60	ND < 0.56
Silver	mg/kg	100	ND <0.10	ND <0.11	ND <0.11	ND <0.12	ND <0.12	ND <0.11	ND <0.11	ND < 0.11	ND <0.10	ND <0.11	ND <0.11	ND <0.12	ND <0.11	ND <0.11	ND <0.11	ND <0.11	ND <0.11	ND <0.10	ND <0.10	ND <0.10	ND <0.10	ND <0.34 ND <0.11	ND <0.12	ND <0.56 ND <0.11
Thallium	mg/kg	8 2500	ND <0.26	ND <0.28	ND <0.28	ND <0.30	ND <0.29	ND <0.28	ND <0.28	ND <0.28	ND <0.26	ND <0.27	ND <0.27	ND <0.29	ND <0.27	ND <0.27	ND <0.28	ND <0.27	ND <0.28	ND <0.25	ND <0.25	ND <0.27	ND <0.27	ND <0.11	ND <0.30	ND < 0.11
Zinc Mercury (Method 7471)	mg/kg	2500	10	13	85	63	27	54	20	26	48	6.8	71	14	14	6.3	4.0	4.2	16	310	84	22	15	48	8.4	28
Cvanide (Method 9012)	mg/kg mg/kg	100	ND <0.021 ND <0.53	ND <0.022 ND <0.56	0.029 ND < 0.56	ND <0.024	0.14 ND <0.50	ND <0.023	0.034	0.051	0.023	ND <0.022	ND <0.021	0.065	ND <0.022	ND <0.022	ND <0.023	ND <0.021	ND <0.022	0.073	ND < 0.023	0.023	ND <0.021	0.032	ND <0.024	0.044
VPH (MADEP Method)	Img/ kg	100	ND <0.33	ND <0.56	ND <0.56	ND <0.59	ND <0.59	ND <0.57	ND <0.57	ND <0.57	ND <0.53	ND <0.54	ND <0.54	ND <0.59	ND <0.55	ND <0.55	ND <0.56	ND <0.53	ND <0.55	ND <0.54	N'D. <0.58	NID < 0.55	ND <0.54	ND <0.54	ND <0.6	ND <0.56
C5-C8 Aliphatics	µg/kg	100000	ND <6200	ND <6500	ND <6900	ND <8800	ND <7900	ND <7000	ND <7000	ND <6900	-5000	277 55500		-7000				,								
C9-C12 Aliphatics	μg/kg	100000	ND <6200	ND <6500	ND <6900	ND <8800	ND <7900	ND < 7000	ND <7000 ND <7000	ND <6900 ND <6900	ND <5900 ND <5900	ND <6500 ND <6500	ND <5900 ND <5900	NID <7800 NID <7800	ND <6400	ND <6400	ND <7600	ND <6000	ND <6100	ND <6400	ND <7600	ND <6400		ND <12000	ND <8000	ND <6800
C9-C10 Aromatics	ug/kg	100000	ND <6200	ND <6500	ND <6900	ND <8800	ND <7900		ND <7000	ND <6900 ND <6900	ND <5900 ND <5900	ND <6500 ND <6500	ND <5900 ND <5900	ND <7800 ND <7800	ND <6400 ND <6400	ND <6400 ND <6400	ND <7600	ND <6000	ND <6100	14000	ND <7600	ND <6400		ND <12000	ND <8000	ND <6800
EPH (MADEP Method)				·		,		112	1417 -1000	140 40,00	IND NOO	IND NOOU	ND >3700	ND \1000	ND <0400	ND <0400	NID <7600	ND <6000	ND <6100	ND <6400	ND <7600	ND <6400	ND <6100	ND <12000	ND <8000	ND <6800
C9-C18 Aliphatics	μg/kg	1000000	ND <10000	ND <10000	ND <11000	ND <11000	ND <11000	ND <11000	ND <11000	ND <11000	ND <10000	ND <11000	ND <11000	ND <12000	ND <10000	ND <10000	ND <11000	ND <11000	ND <10000	ND <11000	ND <11000	ND <10000	ND <10000	ND <11000	277 211000	
C19-C36 Aliphatics	μg/kg	2500000	40000	14000	ND <11000	ND <11000	ND <11000	ND <11000	ND <11000			ND <11000													ND <11000 ND <11000	ND <11000 _
								-1211000	11000	11000	ND < 10000	ND -11000	24000	ND <12000	ND <10000	38000	MD <11000 1	ND <11000	ND <10000	~000 C11000	~!! <11000	NID <10000	~ (1 (X X X X)			
C11-C22 Aromatics	μg/kg	200000	100000	16000	17000	32000	110000	20000	40000	61000		ND <11000	92000	ND <12000 ND <12000	ND <10000 52000	38000 81000			ND <10000	ND <11000 ND <11000	ND <11000 ND <11000	ND <10000 25000	ND <10000 ND <10000			
VOCs (Method 8260)				16000	17000	32000		20000	40000	61000	33000	ND <11000	92000	ND <12000	52000	81000	ND <11000								ND <11000	28000
VOCs (Method 8260) Benzene	μg/kg	2000	ND <5.3	16000 ND <5.6	17000 ND <5.1	32000 ND <5.8	ND <5.3	20000 ND <5.5	40000 ND <350	61000 ND <340	33000 ND <5.0	ND <11000 ND <5.3	92000 ND <300	ND <12000 ND <5.8	52000 ND <5.2	81000 ND <5.4	ND <11000 ND <5.2									
VOCs (Method 8260) Benzene Ethylbenzene	µg/kg µg/kg	2000 80000	ND <5.3 ND <5.3	16000 ND <5.6 ND <5.6	17000 ND <5.1 ND <5.1	32000 ND <5.8 ND <5.8	ND <5.3 ND <5.3	20000 ND <5.5 ND <5.5	40000 ND <350 ND <350	61000 ND <340 ND <340	33000 ND <5.0 ND <5.0	ND <11000 ND <5.3 ND <5.3	92000 ND <300 ND <300	ND <12000 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2	81000 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6	ND <11000 ND <5.6 ND <5.6	25000 ND <270 ND <270	ND <10000 ND <5.4 ND <5.4	18000 ND <5.0 ND <5.0	ND <11000	28000
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether	μg/kg μg/kg μg/kg	2000	ND <5.3 ND <5.3 ND <5.3	16000 ND <5.6 ND <5.6 ND <5.6	17000 ND <5.1 ND <5.1 ND <5.1	32000 ND <5.8 ND <5.8 ND <5.8	ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5 ND <5.5 ND <5.5	40000 ND <350 ND <350 ND <70	61000 ND <340 ND <340 ND <69	33000 ND <5.0 ND <5.0 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59	ND <12000 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2	81000 ND <5.4 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6 ND <5.6 ND <5.6	25000 ND <270 ND <270 ND <64	ND <10000 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0 ND <5.0 ND <5.0	ND <11000 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2 ND <5.2 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene	µg/kg µg/kg µg/kg µg/kg	2000 80000 100	ND <5.3 ND <5.3	16000 ND <5.6 ND <5.6	17000 ND <5.1 ND <5.1	32000 ND <5.8 ND <5.8 ND <5.8 10	ND <5.3 ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5 ND <5.5 ND <5.5 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110	ND <340 ND <340 ND <340 ND <69 ND <110	33000 ND <5.0 ND <5.0 ND <5.0 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260	81000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6 ND <5.6 ND <5.6 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0 ND <5.0	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2 ND <5.2 ND <5.2 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene	μg/kg μg/kg μg/kg	2000 80000 100 4000	ND <5.3 ND <5.3 ND <5.3 ND <5.3	16000 ND <5.6 ND <5.6 ND <5.6 ND <5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1	32000 ND <5.8 ND <5.8 ND <5.8	ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5 ND <5.5 ND <5.5 ND <5.5 ND <5.5	MD <350 ND <350 ND <350 ND <70 ND <110 ND <350	ND <340 ND <340 ND <340 ND <69 ND <110 ND <340	33000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2	81000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6 ND <5.6 ND <5.6 ND <5.6 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <6.0 ND <6.0	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene	µg/kg µg/kg µg/kg µg/kg µg/kg	2000 80000 100 4000 30000	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	16000 ND < 5.6 ND < 5.6 ND < 5.6 ND < 5.6 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17	32000 ND <5.8 ND <5.8 ND <5.8 10 ND <5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5 ND <5.5 ND <5.5 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110	ND <340 ND <340 ND <340 ND <69 ND <110	33000 ND <5.0 ND <5.0 ND <5.0 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2	81000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6	ND <11000 ND <5.6 ND <5.6 ND <5.6 ND <5.6 ND <5.6 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <270	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 16 ND <5.0	ND <11000 ND <5.7	28000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene	µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg	2000 80000 100 4000 30000 NR 500000	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0	32000 NID < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <350	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340	33000 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 ND <5.2	81000 ND < 5.4 ND < 5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <270 ND <270 ND <270	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-xylene Tetrachloroethene 1,2,3-Trichloröbenzene	нд/kg нд/kg нд/kg нд/kg нд/kg нд/kg нд/kg нд/kg нд/kg нд/kg	2000 80000 100 4000 30000 NR 500000 1000	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND < 5.1 ND < 5.1 ND < 5.1 ND < 5.1 T 7.0 ND < 5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <350 ND <350 ND <350 T <350 ND <350	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340	33000 ND <5.0	ND <11000 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2	81000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <270 ND <270 ND <270 4300	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 AD <5.4 ND <5.4	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 16 ND <5.0 ND <5.0 40	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2 11
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichlorobenzene 1,1,1-trichloroethane	µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg µg/kg	2000 80000 100 4000 30000 NR 500000 1000 NR 30000	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 17 7.0 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340	33000 ND <5.0	ND <11000 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300 3000	ND <12000 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 130	81000 ND < 5.4 ND < 5.4 120	ND <11000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 13	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 ND <4.9	ND <10000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <270 ND <270 ND <270	ND <10000 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2.3-Trichloroethane 1,1,2-Trichloroethane		2000 80000 100 4000 30000 NR 500000 1000 NR 500000 1000	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 TO <5.1 ND <5.1 ND <5.1 TO ND <5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3	20000 ND < 5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <1100	33000 ND <5.0	ND <11000 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <100 ND <1100 ND <1100 ND <1100	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2	81000 ND < 5.4	ND <11000 ND <5.2 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 ND <4.9 ND <4.9	ND <10000 ND <4.6	ND <11000 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <270 ND <270 ND <270 ND <270 ND <270	ND <10000 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7 ND <5.7	28000 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane Trichloroethane Trichloroethene	Hg/kg	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 100	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND < 5.1 ND < 5.1 ND < 5.1 ND < 5.1 17 7.0 ND < 5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 28 ND <5.3 ND <5.3	20000 ND < 5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 19000	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100	33000 ND <5.0 SD <5.0 ND <5.0 SD <5.0 SD <5.0 SD <5.0 SD <5.0 SD <5.0 SD <5.0	ND <11000 ND <5.3 22 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <110 ND <300 ND <1100	ND <12000 ND <5.8 79	52000 ND < 5.2 ND < 5.2 ND < 5.2 ND < 5.2 260 ND < 5.2 ND < 5.2 130 ND < 5.2 ND < 5.2 130 ND < 5.2 ND < 5.2 130 ND < 5.2 ND < 5.2 150	81000 ND < 5.4 120 ND < 5.4 Consideration of the constant of the constan	ND <11000 ND <5.2 13 ND <5.2 73 ND <5.2 37	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44	ND <10000 ND <4.6	ND <11000 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270	ND <10000 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7	28000 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Vinyl Chloride		2000 80000 100 4000 30000 NR 500000 1000 NR 500000 1000	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 TO <5.1 ND <5.1 ND <5.1 TO ND <5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3	20000 ND < 5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <350 ND <350 ND <350 ND <350 ND <570 ND <570 ND <570 ND <570 19000	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <1100	33000 ND <5.0	ND <11000 ND <5.3 22 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <100 ND <1100 ND <1100 ND <1100	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2	81000 ND < 5.4	ND <11000 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9	ND <10000 ND <4.6	ND <11000 ND <4.6	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270	ND <10000 ND <5.4 AB ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7	28000 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane Trichloroethane Trichloroethene	HE/kg HE/k	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 3000 600	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 To 17 7.0 ND <5.1	32000 NID < 5.8 NID < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 ND <70 ND <70 ND <70 ND <1000	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <1100 ND <1200 ND <1000	33000 ND <5.0 SD <5.0 ND <5.0 SD <5.0 ND <5.0	ND <11000 ND <5.3	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <1100 ND <1000	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 130 ND <5.2	81000 ND < 5.4	ND <11000 ND <5.2 13 ND <5.2 73 ND <5.2 37 ND <5.0	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8	ND <10000 ND <4.6 Z6 ND <4.6	ND <11000 ND <4.6 ND	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <550	ND <10000 ND <5.4	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 AU ND <5.0	ND <11000 ND <5.7	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 6.5
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-sylenes o-Xylene Tetrachloroethene 1,2,3-Trichloröbenzene 1,1,1-trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method)	HE/kg	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 100	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 T 7.0 ND <5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 28 ND <53 ND <53 12 ND <53 ND <53 ND <53 ND <53	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 ND <570 ND <1100	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <310 ND <310 ND <1100	33000 ND <5.0	ND <11000 ND <5.3	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <59 ND <1110 ND <300 ND <300 ND <300 ND <300 ND <300 ND <1100	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2	81000 ND < 5.4 120 ND < 5.4	ND <11000 ND <5.2 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <10	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8 ND <110	ND <10000 ND <4.6 N	ND <11000 ND <4.6 N	ND <11000 ND <5.6	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <550 ND <100	ND <10000 ND <5.4 ND <5.1 ND <5.4 ND <5.1 ND <5.4	18000 ND <5.0	ND <11000 ND <5.7	28000 ND <5.2 11 ND <5.2 ND <5.2 10 ND <5.2 ND <5.1 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene	HE/kg HE/k	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 3000 600	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 TO <5.1 ND <5.1	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3	20000 ND <5.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 ND <1100 ND <110 ND <110	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <1100	33000 ND <5.0	ND <11000 ND <5.3 ND	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300 ND <1100	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 130 ND <5.2 130 ND <5.2 ND <5.2 130 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 On <5.2 ND <5.2 On	81000 ND < 5.4 120 ND < 5.4	ND <11000 ND <5.2	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8 ND <110 ND <110	ND <10000 ND <4.6 ND <4.5 ND <4.6 ND	ND <11000 ND <4.6 N	ND <11000 ND <5.6 N	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 ND <100 ND <100	ND <10000 ND <5.4 ND <5.1	18000 ND <5.0	ND <11000 ND <5.7 ND <11 ND <110 ND <110	28000 ND <5.2 ND <5.1 ND <5.1 ND <5.1 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene	HE/KE	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 1000 000 000 1000 000 000 000 000	ND <5.3 ND <5.3	16000 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 T 7.0 ND <5.1 ND <740	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3 28 ND <5.3 ND <	20000 ND < 5.5 AS ND < 5.5 AS ND < 5.5 AS ND < 5.5 ND < 1.5 ND < 1.5 ND < 1.1 ND < 1.10 ND < 1.10	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 ND <570 ND <1100	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <3100 ND <1100	33000 ND <5.0	ND <11000 ND <5.3 ND	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300 ND <1100	ND <12000 ND <5.8 ND <100	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.0 150 ND <10	81000 ND < 5.4 ND < 5.1 120 ND < 5.4 ND < 5.4 ND < 5.4 ND < 5.4 ND < 5.1	ND <11000 ND <5.2 ND <10 ND <110 ND <110 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <4.9 N	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 N	ND <11000 ND <5.6 N	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <500 ND <500 ND <500 ND <500 ND <100 ND <100 ND <100 200	ND <10000 ND <5.4 ND <5.1 ND <100 ND <100 ND <100 ND <100	18000 ND <5.0 ND <10	ND <11000 ND <5.7 ND <11 ND <110 ND <110 ND <110	28000 ND <5.2 ND <5.1 ND <5.1 ND <5.1 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthylene Anthracene Benzo(a)anthracene	HE/kg	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 300 600 4000 20000 100000 7000	ND <5.3 ND <5.0 ND	16000 ND < 5.6 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 T 7.0 ND <5.1 ND <40 240	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND <	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND	20000 ND <5.5 ND <1.5	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <570 ND <110 ND <110 ND <110 230	61000 ND <340 ND <340 ND <69 ND <110 ND <340 ND <1100	33000 ND <5.0 ND <100 220 ND <100 970	ND <11000 ND <5.3 ND	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300 ND <1100	ND <12000 ND <5.8	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 27 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 130 ND <5.2 ND <5.2 ND <5.2 150 ND <10 200 560 1800	81000 ND < 5.4 120 ND < 5.4 ND < 5.1 ND < 100 ND < 110 ND < 100 ND < 100 ND < 100	ND <11000 ND <5.2 ND <10 ND <110 ND <110 ND <110 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8 ND <110 ND <110 ND <110 ND <110 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 ND <550 ND <100 ND <100 ND <100 200 250	ND <10000 ND <5.4 ND <100 ND <100 ND <100 ND <100 ND <100	18000 ND <5.0 ND <10 ND <110 ND <110 ND <110 ND <110	ND <11000 ND <5.7 ND <10 ND <110 ND <110 ND <110	28000 ND <5.2 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.2
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthylene Anthracene Benzo(a)anthracene Benzo(a)pyrene	HE/kg	2000 80000 100 4000 30000 NR 5000000 1000 NR 30000 1000 1000 10000 4000 20000 100000 100000 20000	ND <5.3 ND <10 ND <100 ND <100 270 390 1100	16000 ND < 5.6 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110 740 240 ND <110 ND <110 ND <110 ND <110 ND <110	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 6.8 ND < 12 ND < 110 ND < 110 ND < 110 340 320 1000 1000	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3 29 ND <1.3 ND <1.0 ND <	20000 ND <5.5 ND <1.5 ND <1.5 ND <1.5 ND <1.1 ND <1.1 ND <1.10 ND <1.10 350 110 200	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <550 ND <350 ND <350 ND <350 ND <570 ND <570 ND <570 ND <1100 ND <110 230 230 200	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <110	33000 ND <5.0 110 ND <100 220 ND <100 970 1100	ND <11000 ND <5.3 ND	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <1100 ND <100 ND	ND <12000 ND <5.8 ND <120 ND <120 ND <120 ND <120 ND <120	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.0 150 ND <10	81000 ND < 5.4 120 ND < 5.4 ND < 5.1 ND < 100 ND < 110 ND < 100 ND < 100 ND < 100	ND <11000 ND <5.2 ND <10 ND <110 ND <110 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 ND <100 ND <100 ND <100 D <100	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <5.1 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <5.1 ND <110 ND <110 490
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes O-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,2-Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthylene Anthracene Benzo(a)pyrene Benzo(b)fluoranthene	HE/kg	2000 80000 100 4000 30000 100 NR 500000 100 NR 30000 100 3000 100 300 100 100 20000 100000 100000 7000	ND <5.3 ND <10 ND <100 ND <100 100 1100 1100 1700	16000 ND < 5.6 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110	32000 ND < 5.8 ND < 5.8 10 ND < 5.8 10 ND < 5.8 ND < 12 ND < 110 ND < 110 340 320 1000 1400	ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 ND <53 28 ND <53 28 ND <53 12 ND <11 12 ND <11 210 ND <110 1300 2000 4400	20000 ND <5.5	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <350 ND <570 ND <570 ND <570 ND <110 ND <110 ND <110 230 200 330 ND <110	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100	33000 ND <5.0	ND <11000 ND <5.3 ND <6.3 ND <6.6 ND <1.0 ND <110 ND	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 36700 ND <110 ND <110 ND <110 ND <110 ND <110 ND <100 S6700 ND <110 ND <110 ND <100 ND	ND <12000 ND <5.8 ND <120	S2000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 130 ND <5.2 ND <5.2 ND <5.2 ND <5.0	81000 ND < 5.4 120 ND < 5.4 ND < 100	ND <11000 ND <5.2 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <4.9 10 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND <4.110 ND <4.110 ND <4.110	ND <11000 ND <5.6 N	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <250 ND <100 ND <100 200 250 1500 2000	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10	ND <11000 ND <5.7 ND <5.1 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <10 ND <110 490 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-sylenes o-Xylene Tetrachloroethene 1,2,3-Trichloröbenzene 1,1,1-trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthylene Acenaphthylene Anthracene Benzo(a)nntracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(b)fluoranthene	HE/kg	2000 80000 100 4000 30000 100 NR 500000 100 3000 100 300 600 4000 20000 100000 7000 2000 NR	ND <5.3 ND <10 ND <100 100 100 1100 1100 1100 1100 1100 1	16000 ND < 5.6 ND < 11 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110	32000 NID < 5.8 NID < 5.8 NID < 5.8 10 NID < 5.8	ND <5.3 12 ND <5.3 ND	20000 ND <5.5 ND <110	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <370 ND <110 S00 330 ND <1100 350 1500 350 170	61000 ND <340 ND <340 ND <340 ND <340 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <110	33000 ND <5.0 ND <100 220 ND <100 970 1100 ND <100 ND <100 ND <100 ND <100 ND <100 ND <100	ND <11000 ND <5.3 ND <6.3 ND <6.6 ND <1.0 ND <110 ND	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100	ND <12000 ND <5.8 ND <100 ND <100 ND <100 ND <120	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.0 ND	81000 ND < 5.4 ND < 5.1 120 ND < 5.4 ND < 100 ND < 110 ND < 100 N	ND <11000 ND <5.2 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 17 ND <4.9 10 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <250 D <550 ND <100 ND <100 200 250 1500 2000 2000	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <10 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <10 ND <110 ND <110 490 ND <110 290
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,1-trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthene Acenaphthene Benzo(a)anthracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(b)fluoranthene Chrysene	Hg/kg	2000 80000 100 4000 30000 NR 500000 1000 NR 30000 100 300 600 4000 20000 100000 7000 2000 7000 NR 7000	ND <5.3 ND <10 ND <100 ND <100 270 390 1200 1100 1700 1700 600 1600	16000 ND < 5.6 ND < 11 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 T 7.0 ND <5.1 ND <10 ND <110	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 12 ND < 110 ND < 110 340 320 1000 1000 1400 540 1200	ND <53 ND <50 ND <110 ND <110 1300 2000 4400 39900 5600 1900 4500	20000 ND <5.5 ND <10 ND <110	40000 ND <350 ND <350 ND <70 ND <110 ND <350 ND <110 230 200 330 ND <110 560 170 640	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <110 S10 ND <110 S90	33000 ND <5.0 ND <5	ND <11000 ND <5.3 ND <5.1 ND <110	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 260 ND <5.2 130 ND <5.2 130 ND <5.2 ND <5.2 130 ND <5.2 150 ND <5.2 150 150 150 1800 2800 2800 2600	81000 ND < 5.4 120 ND < 5.4 ND < 11 ND < 100	ND <11000 ND <5.2 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <9.8 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND <5.10 ND <110	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <250 ND <100 ND <100 D <1	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 16 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <10 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <10 ND <110 490 ND <110 290 ND <110 290 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-sylenes o-Xylene Tetrachloroethene 1,2,3-Trichloröbenzene 1,1,1-trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthylene Acenaphthylene Anthracene Benzo(a)nntracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(b)fluoranthene	HE/kg	2000 80000 100 4000 30000 NR 5000000 1000 NR 30000 1000 1000 10000 10000 100000 100000 NR 300 600	ND <5.3 ND <5.0 ND <5.0 ND <100 ND <100 100 1100 1700 600 600 1600 2200	16000 ND < 5.6 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110 N	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND <	ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 ND <5.3 28 ND <5.3 ND <5.3 28 ND <5.3 ND <5.0 ND <110 1300 2000 5600 1900 4500 4500 8200	20000 ND <5.5	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <350 ND <100 ND <110 ND <110 S00 ND <110 ND <110 S00 ND <110 S00 S00 ND <110 S00 S00 S00 S00 S00 S00 S00 S00 S00	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <110 N	33000 ND <5.0 ND <10 ND <100 220 ND <100 970 1100 ND <100 2300	ND <11000 ND <5.3 ND <5.10 ND <110	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	52000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 150 ND <5.2 150 ND <5.2 150 ND <10 200 560 1800 2800 2800 1100	81000 ND < 5.4 ND <	ND <11000 ND <5.2 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 44 ND <4.9 AH ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND <4.110 ND <110	ND <11000 ND <5.6 ND	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 D <270 ND <270 ND <270 D <270 ND <270 ND <270 D <2000 D <2000 2000	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 16 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <10 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <110 ND <110 490 ND <110 290 ND <110 ND <110 ND <110 290 ND <110 ND <110 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloröbenzene 1,1,1-trichloroethane Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthylene Anthracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(k)fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene	HE/kg	2000 80000 100 4000 30000 NR 500000 100 3000 100 3000 100 3000 100 300 600 400 20000 100000 7000 2000 NR 7000 NR 7000 NR 7000 1000000 400	ND <5.3 ND <10 ND <100 ND <100 ND <100 1700 600 1700 600 1600 2200 ND <100	16000 ND < 5.6 ND < 10 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110	32000 NID < 5.8 NID < 5.8 NID < 5.8 10 NID < 5.8 10 NID < 5.8 NID < 5	ND <5.3 ND <5.	20000 ND <5.5 ND <10 ND <110	40000 ND <350 ND <350 ND <70 ND <110 ND <110 ND <350 ND <550 ND <550 ND <570 ND <570 ND <570 ND <110 380 ND <110 640 380 ND <110	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <110 ND <110 ND <110 ND <110 S50 ND <110	33000 ND <5.0 ND <100 220 ND <100 970 1100 ND <100 2300 2300	ND <11000 ND <5.3 ND <6.6 ND <10 ND <110	92000 ND <300 ND <300 ND <59 ND <110 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	\$2000 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 260 ND <5.2 130 ND <5.2 ND <5.2 130 ND <5.2 ND <5.2 150 ND <5.2 150 ND <10 200 560 1800 2800 2200 2600 1100 2800	81000 ND <5.4 ND <5.1 120 ND <5.4 ND <5.4 ND <5.1 120 ND <100	ND <11000 ND <5.2 13 ND <5.2 73 ND <5.2 73 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 18 ND <4.9 10 ND <1.0	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND <5.10 ND <110	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 D <270 ND <270 D <270	ND <10000 ND <5.4 ND <5.1 ND <100 ND <100	18000 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <110 ND <110 120 ND <110 490 ND <110 290 ND <110 290 ND <110 1300
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthylene Anthracene Benzo(a)anthracene Benzo(b)fluoranthene Benzo(b)fluoranthene Chrysene Fluoranthene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene	HE/kg	2000 80000 100 4000 30000 100 NR 500000 1000 3000 600 4000 20000 100000 7000 2000 NR 7000 100000 100000 7000 100000 7000 70	ND <5.3 ND <6.3 ND <5.3 ND <100 ND <100 270 390 1200 1100 1100 1700 600 1600 2200 ND <100	16000 ND < 5.6 ND < 11 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 TO ND <5.1 17 7.0 ND <5.1 ND <10 ND <110	32000 NID < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 12 ND < 110 ND < 110 ND < 110 340 1400 540 1200 2100 ND < 110 ND < 110 ND < 1110 ND < 1110	ND <5.3 ND <5.	20000 ND < 5.5 ND < 110	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <310 ND <350 ND <350 ND <350 ND <350 ND <350 ND <350 ND <370 ND <110 S80 ND <100 S80 ND <110 S90 ND <110	61000 ND <340 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <110	33000 ND <5.0 Self	ND <11000 ND <5.3 ND <10 ND <110	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 130 ND <5.2 ND <5.2 ND <5.2 150 ND <5.2 150 ND <10 200 200 200 2800 2600 1100 2800 2800 6700	81000 ND < 5.4 ND < 100 ND < 110 ND < 100 Add Add ND < 100	ND <11000 ND <5.2 13 ND <5.2 13 ND <5.2 16 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 18 ND <4.9 17 ND <4.9 10 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND <4.10 ND <110	ND <11000 ND <5.6 ND <5.10 ND <110	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <200 2000 2000 2000 2000 2000 2000 360 120 1300 ND <100	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <110 ND <110 490 ND <110 290 ND <110 ND <110 ND <110 290 ND <110 ND <110 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloröbenzene 1,1,1-trichloroethane Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthylene Anthracene Benzo(a)pyrene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(k)fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene Fluoranthene	Hg/kg Hg/k	2000 80000 100 4000 30000 NR 30000 1000 NR 30000 1000 300 600 4000 20000 100000 7000 2000 7000 NR 7000 1000000 400 7000 1000000	ND <5.3 ND <10 ND <100 ND <100 ND <100 270 390 1200 1100 1700 1700 1600 2200 ND <100 ND <100 990	16000 ND < 5.6 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110 N	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 12 ND < 110 ND < 110 ND < 110 340 320 1000 1400 540 1200 2100 ND < 110	ND <53 ND <510 ND <110 ND <110 ND <110 ND <100 ND	20000 ND <5.5 ND <5.1 ND <110	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <350 ND <350 ND <570 ND <570 ND <570 ND <570 ND <570 ND <570 ND <110 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <110 D <110 D <110 D <110 D <110 D <110 ND <110	33000 ND <5.0 ND <100 ND <100 ND <100 ND <100 1200 2300 230 ND <100 2300	ND <11000 ND <5.3 ND <5.1 ND <110	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	\$2000 ND <5.2 ND <5.2 ND <5.2 SECOND <5.	81000 ND <5.4	ND <11000 ND <5.2 13 ND <5.2 73 ND <10 ND <110	ND <11000 ND <4.9 18 ND <4.9 17 ND <4.9 18 ND <4.9 17 ND <4.9 10 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND	ND <11000 ND <5.6 ND <110	25000 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 ND <100 200 250 1500 2000 2000 860 120 1300 ND <100 ND <100 ND <100 260	ND <10000 ND <5.4 ND <100	18000 ND <5.0 ND <5.0 ND <5.0 ND <5.0 ND <5.0 16 ND <5.0 ND <10 ND <110	ND <11000 ND <5.7 ND <110	28000 ND <5.2 11 10 ND <5.2 ND <5.2 ND <5.2 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.1 ND <110 ND <110 ND <110 490 ND <110 490 ND <110 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichlorobenzene 1,1,1-trichloroethane Trichloroethane Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthylene Anthracene Benzo(a)phrene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(b)fluoranthene Benzo(c)rysene Fluoranthene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene Phenanthrene Pyrene	HE/kg HE/K	2000 80000 100 4000 30000 100 NR 500000 100 100 3000 100 3000 100 300 600 4000 100000 100000 NR 7000 NR 7000 100000 400 7000 1000000 1000000 1000000 1000000 1000000	ND <5.3 ND <5.0 ND <100 ND <100 ND <100 1700 1700 600 1100 1700 600 1000 2200 ND <100 ND <100 ND <100 1700 600 1000 900 2200 ND <100 ND <100 ND <100 ND <100 000 1000 000 000 000 000 000 000 00	16000 ND < 5.6 ND < 5.6	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 T 7.0 ND <5.1 ND <10 ND <110	32000 ND <5.8 ND <5.8 ND <5.8 10 ND <5.8 N	ND <5.3 ND <5.	20000 ND <5.5 ND <5.10 ND <110	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <350 ND <370 ND <110 ND <110 ND <110 ND <110 ND <110 S60 170 640 380 ND <110	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <1100 ND <110 ND	33000 ND <5.0 ND <100 ND <100 P70 1100 ND <100 1200 2300 2300 2300 2300 2000	ND <11000 ND <5.3 ND <5.1 ND <110	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110 ND <10 990	ND <12000 ND <5.8 N	52000 ND <5.2 ND <5.2 ND <5.2 ND <5.2 260 ND <5.2 ND <5.2 ND <5.2 ND <5.2 ND <5.2 130 ND <5.2 ND <5.2 ND <5.2 150 150 1800 2800 2800 2600 1100 2800 6700 6300 ND <100 7600 6600	81000 ND <5.4 ND <100	ND <11000 ND <5.2 13 ND <5.2 13 ND <5.2 13 ND <5.2 13 ND <10 ND <110	ND <11000 ND <4.9 ND <4.9 ND <4.9 ND <4.9 2300 ND <4.9 ND <4.9 18 ND <4.9 17 ND <4.9 18 ND <4.9 10 ND <10 ND <110	ND <10000 ND <4.6 ND	ND <11000 ND <4.6 ND <4.10 ND <110	ND <11000 ND <5.6 ND <110	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 200 250 1500 2000 2000 860 120 1300 ND <100 260 260 1600	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10 ND <110 ND <110	ND <11000 ND <5.7 ND <10 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 11 ND <5.2 ND <10 ND <110 ND <110 490 ND <110
VOCs (Method 8260) Benzene Ethylbenzene Methyl tert butyl ether Naphthalene Toluene M/P-xylenes o-Xylene Tetrachloroethene 1,2,3-Trichloroethane 1,1,1-trichloroethane 1,1,2-Trichloroethane Trichloroethene Vinyl Chloride PAHs (MADEP Method) 2-Methylnaphthalene Acenaphthene Acenaphthylene Anthracene Benzo(a)anthracene Benzo(b)fluoranthene Benzo(b)fluoranthene Chrysene Fluoranthene Fluoranthene Fluorene Indeno(1,2,3-cd)pyrene	Hg/kg Hg/k	2000 80000 100 4000 30000 NR 30000 1000 NR 30000 1000 300 600 4000 20000 100000 7000 2000 7000 NR 7000 1000000 400 7000 1000000	ND <5.3 ND <10 ND <100 ND <100 ND <100 270 390 1200 1100 1700 1700 1600 2200 ND <100 ND <100 990	16000 ND < 5.6 ND < 100	17000 ND <5.1 ND <5.1 ND <5.1 ND <5.1 ND <5.1 17 7.0 ND <5.1 ND <10 ND <110 N	32000 ND < 5.8 ND < 5.8 ND < 5.8 10 ND < 5.8 ND < 12 ND < 110 ND < 110 ND < 110 340 320 1000 1400 540 1200 2100 ND < 110	ND <53 ND <510 ND <110 ND <110 ND <110 ND <100 ND	20000 ND <5.5 ND <5.1 ND <110	40000 ND <350 ND <350 ND <350 ND <70 ND <110 ND <350 ND <350 ND <350 ND <350 ND <570 ND <570 ND <570 ND <570 ND <570 ND <570 ND <110 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	61000 ND <340 ND <340 ND <340 ND <69 ND <110 ND <340 ND <340 ND <340 ND <340 ND <340 ND <1100 ND <110 D <110 D <110 D <110 D <110 D <110 ND <110	33000 ND <5.0 ND <100 ND <100 ND <100 ND <100 1200 2300 230 ND <100 2300	ND <11000 ND <5.3 ND <5.1 ND <110	92000 ND <300 ND <300 ND <300 ND <59 ND <110 ND <300 ND <300 ND <300 ND <300 ND <1100 ND <110	ND <12000 ND <5.8 ND <120	\$2000 ND <5.2 ND <5.2 ND <5.2 SECOND <5.	81000 ND <5.4	ND <11000 ND <5.2 13 ND <5.2 73 ND <10 ND <110	ND <11000 ND <4.9 18 ND <4.9 17 ND <4.9 18 ND <4.9 17 ND <4.9 10 ND <110	ND <10000 ND <4.6 ND <4.0 ND <4.00 N	ND <11000 ND <4.6 ND <4.10 ND <110	ND <11000 ND <5.6 ND <5.10 ND <110	25000 ND <270 ND <270 ND <270 ND <64 ND <100 ND <270 ND <100 ND <100 200 250 1500 2000 2000 860 120 1300 ND <100 ND <100 2000 2000 860 120 1300 ND <100 2000	ND <10000 ND <5.4 ND <5.1 ND <100	18000 ND <5.0 ND <10 ND <110 ND <110	ND <11000 ND <5.7 ND <110 ND <110	28000 ND <5.2 11 ND <5.2 ND <5.2 ND <5.2 11 ND <10 ND <110 120 ND <110 490 ND <110 290 ND <110 940

---- Not applicable
ND <X: Compound not detected above laboratory reporting limit
VPH: Volatile petroleum hydrocarbons
EPH: Extractable petroleum hydrocarbons
PAHs: Polycyclic aromatic hydrocarbons
PCBs: Polychlorinated biphenyls
VOCs: Volatile organic compounds
RC GW-1: MADEP Reportable concentrations for GW-1 groundwater areas.
RC GW-2: MADEP Reportable concentrations for GW-2 groundwater areas.
Color-shaded values indicate exceedence of reportable concentration.

SAH LCK



Table 2 Summary of Groundwater Analytical Data and Objectives

Former Nu-StyleCompany, Inc. Grove Street Franklin, Massachusetts

Phase II Comprehensive Site Assessment Report Prepared for the County of Norfolk, Massachusetts

January 2007

			MW-01	MW-01	MW-02	MW-03	MW-04	MW-05
			841061208-27	841061208-28	841061208-30	841061208-32	841061208-29	841061208-31
		MCP	12/8/2006	12/8/2006	12/8/2006	12/8/2006	12/8/2006	12/8/2006
		RCGW-2	Primary	Duplicate 1	Primary	Primary	Primary	Primary
Groundwater parameters	UNITS							
pH	S		5.97	5.97	6.59	6.33	5.97	NA
Specific Conductance	uMhos/cm		464	464	1727	1534	2010	NA
Temperature	C deg		13.1	13.1	8.3	7.6	10.5	NA
Turbidity	ntu		36	36	50	500	15	NA
Dissolved Oxygen	mg/l		0.4	0.4	7.7	6.9	2.6	NA
ORP	mv		-35.0	-35.0	59	93.2	45.1	NA
Metals (Method 6010)								
Barium	mg/l	50	0.042	0.038	0.15	0.21	0.14	0.83
Beryllium	mg/l	0.05	ND < 0.0010	ND < 0.0010	ND < 0.0010	0.0087	ND < 0.0010	0.0018
Cadmium	mg/l	0.004	ND < 0.0020	0.0034				
Chromium	mg/l	2	ND < 0.010	ND < 0.010	ND < 0.010	0.036	ND < 0.010	0.092
Copper	mg/l	100	ND < 0.010	ND < 0.010	0.015	0.018	ND < 0.010	0.073
Lead	mg/l	0.01	0.014	0.012	ND < 0.0040	0.098	ND < 0.0040	1.9
Nickel	mg/l	0.2	ND < 0.010	ND < 0.010	0.15	0.054	0.017	0.12
Zinc	mg/l	0.9	0.023	0.015	0.057	0.17	0.028	0.73
VOCs (Method 8260)								
Methyl tert butyl ether (MTBE)	μg/l	1000	ND <1.0	ND <1.0	ND <1.0	ND <1.0	2.1	ND <1.0
cis-1,2-Dichloroethene	μg/l	100	ND <1.0	ND <1.0	ND <1.0	ND <1.0	8	ND <1.0
Tetrachloroethene	μg/l	50	ND <1.0	ND <1.0	6.6	43	240	ND <1.0
1,1,1-trichloroethane	μg/l	4000	ND <1.0	ND <1.0	ND <1.0	ND <1.0	1.8	ND <1.0
Trichloroethene	μg/I	30	ND <1.0	ND <1.0	6.6	40	150	ND <1.0

--- Not applicable

ND <X: Compound not detected above laboratory reporting limit

NA: Not analyzed

VOCs: Volatile organic compounds

RC GW-1: MADEP Reportable concentrations for GW-1 groundwater areas. RC GW-2: MADEP Reportable concentrations for GW-2 groundwater areas. Color-shaded values indicate exceedence of reportable concentration.



Table 3 Groundwater Elevation Measurements for On-Site Monitoring Wells Measured December 8, 2006

Former Nu-Style Company, Inc. Facility 87 Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Report Prepared for Norfolk County, Massachusetts

January 2007

Location	Time	Depth to Water (feet from PVC)	Absolute Elevation of PVC ^a (feet)	Groundwater Elevation (feet)
MW-1	0943	4.16	100.35	96.19
MW-2	1300	7.96	98.54	90.58
MW-3	1412	8.18	99.73	91.55
MW-4	1130	7.56	98.23	90.67
MW-5	1515	8.49	104.47	95.98

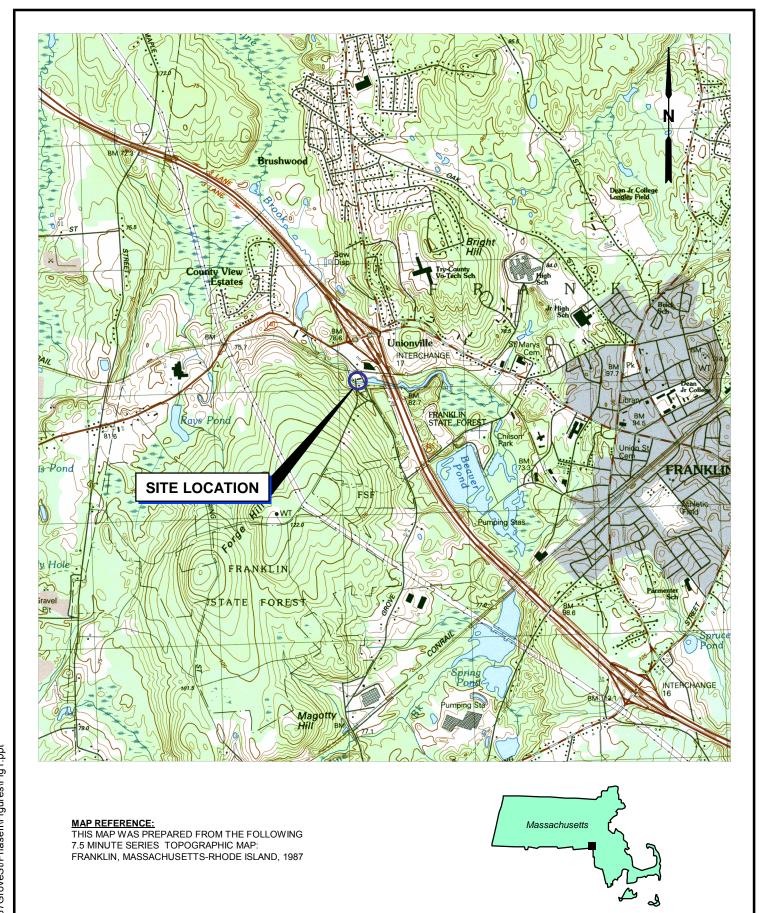
^aelevation data from survey conducted December 4, 2006

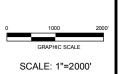
Created by <u>SAH</u> Reviewed by <u>LCK</u>



FIGURES

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT







FOUNDRY CORPORATE OFFICE CENTER
275 PROMENADE ST, SUITE 350, PROVIDENCE RI 02908
401-861-3070 www.FandO

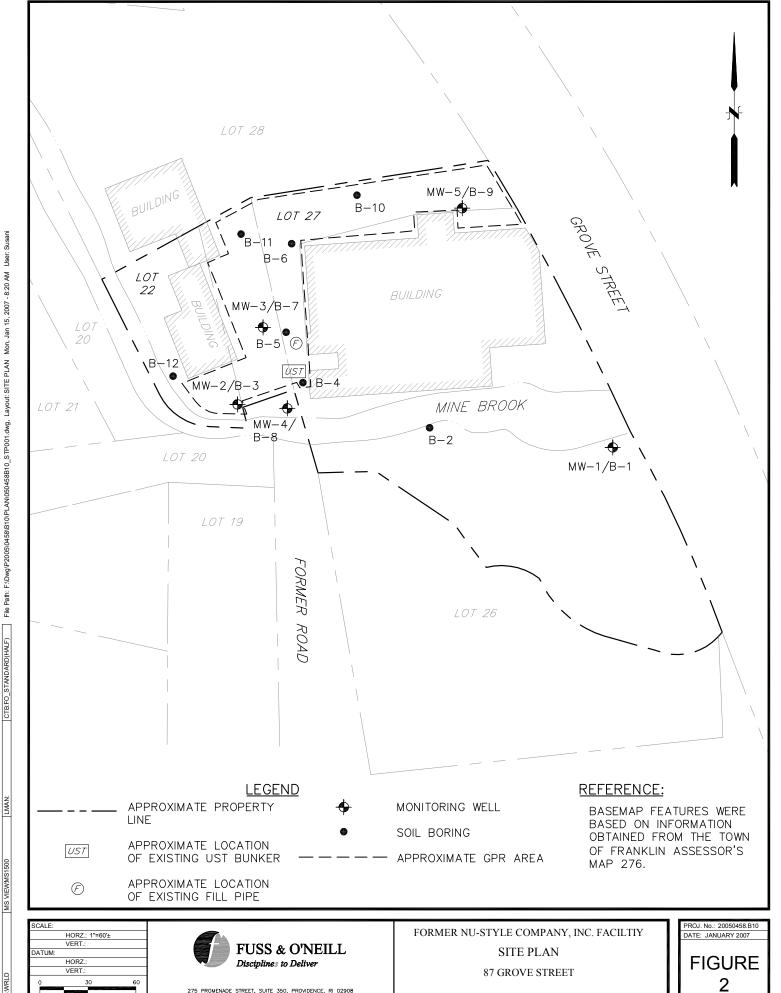
NORFOLK COUNTY, MASSACHUSETTS
SITE LOCATION MAP

FORMER NU-STYLE COMPANY, INC.

87 GROVE STREET FRANKLIN, MASS.

PROJ. No: 20050458.B10 DATE: JANUARY 2007

FIGURE 1



MASSACHUSETTS

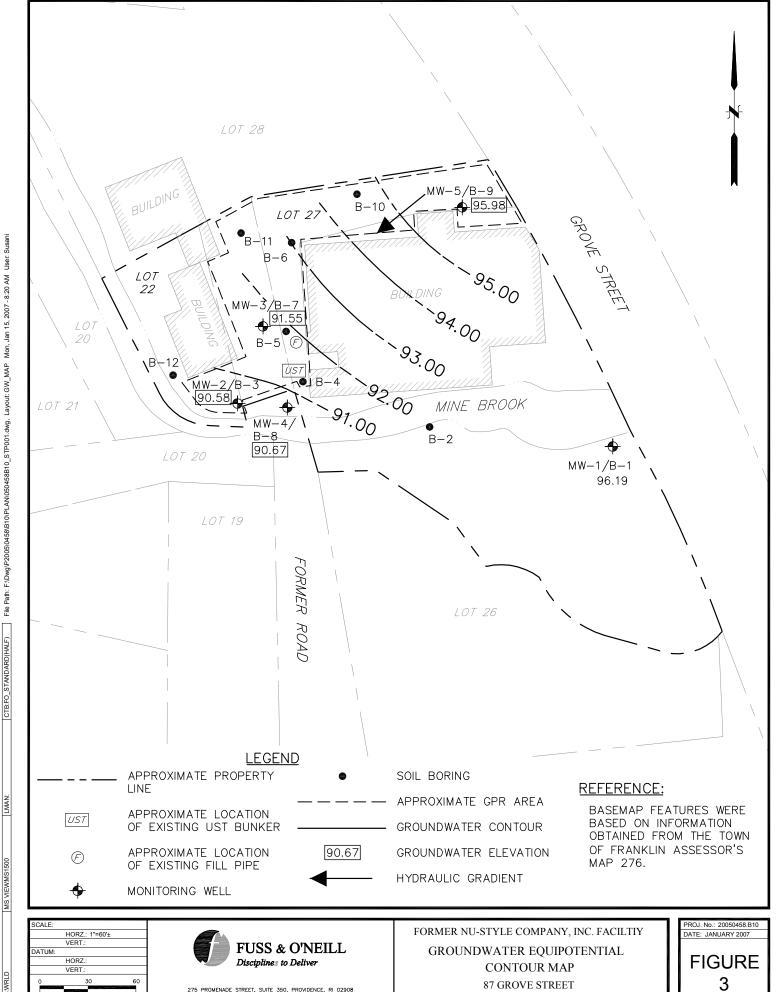
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GRAPHIC SCALE

401.861.3070



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APPENDIX A

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT

SOIL BORING LOGS AND MONITORING WELL COMPLETION REPORTS

Township/Range: Franklin, Massachusetts

Site Id: MW-01

Project Number: 2005-0458 B10



Location: See map

Description: Monitoring Well, Shallow Date(s): 11/30/06 - 11/30/06

Completed Depth: 12.00'

Total Depth: 12.00'

Remarks: Field Instrument: Photovac 2020

Datum: Assumed

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Logged By: S. Hubbs

Driller: S. Perry Contractor: New England Geotech Borehole Dia.: 2.00in

Drilling Method: Geoprobe

Blank Casing: type: PVC

dia: 2.00in

fm: 0.0'

to: 2.00'

Screens: type: Slotted size: 0.010in dia: 2.00in fm: 2.00' to: 12.00'

Annular Fill:

type: Concrete to: 0.50' fm: 0.00' type: Bentonite Grout fm: 0.50' to: 1.00' type: #2 Sand fm: 1.00' to: 12.00'

type: fm: to:

Space Spac					type:	fm:	to:	
-01	Elevation	Depth		Recovery	Material Description		РНОТ	
-10 - 10 -	-2- -4- -6- -8- -10- -12- -14- -16-	2 - 4 - 6 - 8 - 10 - 12 - 14 - 16 - 16 - 16 - 16 - 16 - 16 - 16	N/A -01	R		AS		O ppm

Township/Range: Franklin, Massachusetts

Unable to penetrate after several tries, cannot

move location due to nearby inclined surface and

Site Id: MW-02

Project Number: 2005-0458 B10



Location: See map Description: Monitoring Well, Shallow

Refusal at 12 feet.

drain/water line.

Datum: Assumed Ground Elevation: 0.00' Logged By: S. Hubbs Contractor: New England Geotech

dia: 2.00in

Driller: S. Perry Borehole Dia.: 2.00in

Date(s): 11/30/06 - 11/30/06 Completed Depth: 12.00'

Coordinate X: 0.00

Drilling Method: Geoprobe

to: 2.00'

Total Depth: 12.00' Remarks: Field Instrument: Photovac 2020

Coordinate Y: 0.00

Blank Casing: type: PVC

Screens: type: Slotted size: 0.010in dia: 2.00in fm: 2.00'

fm: 0.0'

to: 12.00'

Annular Fill:

type: Concrete to: 0.50' fm: 0.00' type: Bentonite Grout fm: 0.50' to: 1.50' type: #2 Sand fm: 1.50' to: 12.00' type: fm· to.

				type: type:		fm: fm:	to: to:	
Elevation	Depth	Sample No.	Recovery	Material Description	g	Well Construction MP. EL. 0.00	РНОТ	
-	-	-05		0-0.5': Sand, F-M and gravel; unburned coal/ash; light olive gray (5Y 5/2), dry. 0.5-0.7': CONCRETE. 0.7-5.0': Sand, F-M and gravel; unburned coal/ash; dusky brown (5YR 2/2), moist.		SP CR	4	U ppm
-2-	2	N/A	<u> </u>					
-4-	4	-06		Sand, F-M and gravel; moderate brown (5YR 4/4), moist to wet at 6.0 feet.				0 ppm
-6 -	6	N/A				SP		
-8-	8							
-10-	10 -							
-12-	12-			Boulder or rock. End of boring at 12 feet.		BD		
-14-	14 —							
-16-	16 –							
Chec	cked B	 By:					Page 1 of	1

Township/Range: Franklin, Massachusetts

Site Id: MW-03

Project Number: 2005-0458 B10



Location: See map Description: Monitoring Well, Shallow

Date(s): 11/30/06 - 11/30/06 Completed Depth: 10.50'

Total Depth: 10.50'

Remarks: Field Instrument: Photovac 2020

Refusal at 10.5 feet.

Datum: Assumed

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Logged By: S. Hubbs Contractor: New England Geotech

Driller: S. Perry Borehole Dia.: 2.00in

Drilling Method: Geoprobe

Blank Casing: type: PVC

dia: 2.00in fm: 0.0' to: 2.50'

Screens: type: Slotted size: 0.010in dia: 2.00in fm: 2.50' to: 10.50'

Annular Fill:

type: Concrete fm: 0.00' to: 0.50' type: Bentonite Grout fm: 0.50' to: 1.50' type: #2 Sand fm: 1.50' to: 10.50' tvpe: fm· to.

			fm: to: fm: to:				
Elevation Depth Sample No.	Recovery	Material Description	Graphic Log	USCS Code	Well Construction MP. EL. 0.00	РНОТ	
-13 -2 - 2 - N/A -4 - 4		0-1.0': SAND, F-M; some gravel; dusky brown (5YR 2/2), moist. 1.0-5.0': SAND, F; trace gravel; dusky yellow (5Y 6/4). 5.0-6.1': SAND, F-M, moderate brown (5YR 4/4) to black (N1) at 5.8 feet, moist to wet at 5.8 feet. 6.1-10': SAND, F-M; some gravel; moderate brown (5YR 4/4), wet. Sand, F-M and gravel; dusky yellow (5Y 6/4), wet. Boulders or bedrock. End of boring at 10.5 feet.		SW SP	Page 1 of	O ppm O ppm	

Township/Range: Franklin, Massachusetts

Site Id: MW-04

Project Number: 2005-0458 B10



Location: See map

Description: Monitoring Well, Shallow Date(s): 11/30/06 - 11/30/06

Completed Depth: 10.50' Total Depth: 10.50'

Remarks: Field Instrument: Photovac 2020

Refusal at 10.5 feet.

Datum: Assumed

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Logged By: S. Hubbs

Contractor: New England Geotech

Driller: S. Perry Borehole Dia.: 2.00in

Drilling Method: Geoprobe

Blank Casing: type: PVC

dia: 2.00in

fm: 0.0'

to: 2.50'

Screens: type: Slotted size: 0.010in dia: 2.00in fm: 2.50' to: 10.50'

Annular Fill:

type: Concrete to: 0.50' fm: 0.00' type: Bentonite Grout fm: 0.50' to: 1.50' type: #2 Sand to: 10.50' fm: 1.50'

type: fm: type: fm:	to:
Material Description Order of the control of the c	Construction EL. 0.00
0-1.1': Sand, F and silt; trace brick and coal; dusky brown (5YR 2/2), dry. 1.1-5.0': SAND, M; trace garvel; dusky yellow (5Y 6/4), moist. SM -2- 2- N/A -4- 4-	U ppm
SAND, F-M; trace gravel; 1/4 inch black discrete band at 5.6 feet; dusky yellow (5Y 6/4) with oxidized orange at 6.8 feet; moist to wet at 6.8 feet. SP N/A -8 - 8 -	0 ppm
-10 - 10 - N/A Sand, M-C and gravel; moderate olive brown (5Y 4/4), wet. End of boring at 10.5 feet. -12 - 12	O ppm
Checked By:	Page 1 of 1

Township/Range: Franklin, Massachusetts

Site Id: MW-05

Project Number: 2005-0458 B10



Driller: S. Perry

to: 2.00'

Location: See map

Description: Monitoring Well, Shallow Date(s): 12/01/06 - 12/01/06

Completed Depth: 12.00'

Total Depth: 12.00'

Remarks: Field Instrument: Photovac 2020

Datum: Assumed

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Logged By: S. Hubbs

Contractor: New England Geotech Borehole Dia.: 2.00in

Drilling Method: Geoprobe

Blank Casing: type: PVC

dia: 2.00in fm: 0.0'

Screens: type: Slotted size: 0.010in dia: 2.00in fm: 2.00' to: 12.00'

Annular Fill:

to: 0.50' type: Concrete fm: 0.00' type: Bentonite Grout fm: 0.50' to: 1.50' type: #2 Sand fm: 1.50' to: 12.00'

Material Description Same and coal pieces					type: type:		fm: fm:	to: to:	
10	Elevation	Depth	Sample No.	Recovery		Graphic Log			
-2- 2-N/A -4- 47-20 -6- 67-10- 10-N/A -8- 812- 1214- 1416- 1614- 1416- 1617-1218- 1618- 1819- 18	_	-	1 '		I drv. Loose.	0	SW	\$ 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0 0	U ppm
-6- 6-	_	-	N/A			0 0			
-8- 8- -10- 10- N/A -10- 10- N/A -12- 12- -14- 14- -16- 16- -16- 16-	-6-	6 —			5.0-6.0': Same as above. 6.0-10': SILT, clayey, light olive gray (5Y 5/2), wet.	0 0			0 ppm
11.2-12: SILT, clayey, light olive gray (57 5/2), wet. SW SW ML End of boring at 12 feet. -14- 14- 14	-8 -						ML		
-12 - 12 - End of boring at 12 feet. -14 - 14 -	-10 -	10 -	 N/A 		10-11.2': Sand, F-M and gravel; duksy yellow (5Y 6/4), wet. 11.2-12': SILT, clayey, light olive gray (5Y 5/2), wet.		J#		0 ppm
	-12-	12 -		<u></u>	End of boring at 12 feet.	= = = =		[6594 1555H	
	-14 -	14							
Checked By: Page 1 of 1	_	-	-						
	Chec	ked B	By:					Page 1 of	1

Project Location: Franklin, Massachusetts

Site Id: B-02

Project Number: 2005-0458 B10

Back Fill:



Driller: S. Perry

Location: See map Datum: Logged By: S. Hubbs Description: Soil Boring Ground Elevation: 0.00'

Contractor: New England Geotech Borehole Dia.: 2.00in

Date(s): 11/30/06 - 11/30/06 Coordinate X: 0.00 Drilling Method: Geoprobe

Total Depth: 7.50' Coordinate Y: 0.00

Remarks: Field Instrument: Photovac 2020

Refusal at 7.5 feet.

type: Native Material fm: 0.00' to: 7.50' type: fm: to: fm:

	type: type: type:	fm: fm: fm:	to: to: to:	
Elevation Depth Sample No. Recovery	Material Description	Graphic Log	USCS Code	РНОТ
	0-0.2': Sand, F and silt; some organics; leaf litter; dusky brown (5YR 2/2), dry. 0.2-5.0': Sand, F-M and gravel; dark yellowish brown (10YR 4/2), moist.		SM	U ppm
$\begin{bmatrix} -2 - & 2 - \\ - & - \end{bmatrix}_{N/A}$				
-4- 4	5.0-5.3': BRICK, red. 5.3-7.5': Sand, F-M and gravel; oxidized from 5.3 to 5.8 feet; dark yellowish brown (10YR 4/2), wet.		SP	0 ppm
-6- 6- - N/A	(10YR 4/2), wet.		RK	
-8- 8-	Rock or brick. End of boring at 7.5 feet.			
-10 - 10 -				
 -12 				
-14 - 14 -				
-16 - 16 - 				
-18 - 18 - Checked By:		Page 1	of 1	

Checked By:

Project Location: Franklin, Massachusetts

Site Id: B-04

Project Number: 2005-0458 B10



Page 1 of 1

NADE STREET, SUITE 350, PROVIDENCE, RI 02908 Location: See map Datum: Logged By: S. Hubbs Driller: S. Perry Description: Soil Boring Contractor: New England Geotech Borehole Dia.: 2.00in Ground Elevation: 0.00' Date(s): 11/30/06 - 11/30/06 Coordinate X: 0.00 Drilling Method: Geoprobe Coordinate Y: 0.00 Total Depth: 9.00' Back Fill: fm: 0.00' Remarks: Field Instrument: Photovac 2020 type: Native Material to: 9.00' type: fm: to: Refusal at 9.0 feet. Pulled piece of granite bridge abutment out of type: fm: to: fm: to: abandoned first hole. type: fm: type: to: 2 ġ Code Elevation Recovery Sample Graphic Material Description **NSCS** Depth PHOT U ppm -07 0 SW 0-0.3': Sand, F-M and gravel; light olive gray (5Y 2/2), dry. 0.3-5.0': SAND, F-M; some gravel; dusky brown (5YR 2/2). $2 \dashv_{N/A}$ -2 4 SP -08 0 ppm SAND, F-M; some gravel; dusky brown (5YR 2/2), wet at 6.0 feet. -6 6 -ĦΝ/A -8 8 RK Brick or rock. End of boring at 9.0 feet. 10 -10 - -12 - 12 –14· 14 -16 16 -18 18

Date(s): 11/30/06 - 11/30/06

Project Location: Franklin, Massachusetts

Site Id: B-05

Project Number: 2005-0458 B10



Location: See map Description: Soil Boring Datum:

Logged By: S. Hubbs Driller: S. Perry Contractor: New England Geotech Borehole Dia.: 2.00in

Coordinate X: 0.00 Drilling Method: Geoprobe

Total Depth: 9.00'

Coordinate Y: 0.00

Ground Elevation: 0.00'

Remarks: Field Instrument: Photovac 2020

Refusal at 9.0 feet.

Back Fill: type: Native Material fm: 0.00'

to: 9.00' type: fm: to: type: fm: to:

Section Page Page					type: fm type: fm	:	to: to:	
0 ppm DA-14-Y Sond, F-M and grovel; trace coal/ash; dark yellowish brown (10YR 4/2), moist. SW SW SW SW SW SW SW S				Recovery	Material Description	Graphic Log		РНОТ
-4 - 4 -	_	-				0	Į.	0 ppm
-6- 6- 8- N/A -8- 8-	-2-	2	N/A					
5.0-6.2: Same as above. 6.2-9.0: Sand, M and gravel: dusky yellow (SY 6/4), wet. -8-8-8- End of boring at 9.0 feet. -12-1214-1416-1616-16-	-4-	4 —					SP	
-8 - 8 -	-6-	6 —		/	5.0-6.2': Same as above. 6.2-9.0': Sand, M and gravel; dusky yellow (5Y 6/4), wet.			0 ppm
End of boring at 9.0 feet. -10 - 10	_8 _						SW	
-12- 12- -14- 14- -16- 16- - 1	-0	-			End of boring at 9.0 feet.			
-14 - 14	-10-	10 -						
-16- 16- 	-12-	12 —						
	_ _14 _	14 —						
	-16-	- 16 <i>-</i> -						
		-						
-18 - 18 - Checked By: Page 1 of 1	-18- Checke	18 – ed By:				Page 1	of 1	

Project Location: Franklin, Massachusetts

Site Id: B-06

Project Number: 2005-0458 B10



Location: See map

Datum:

Driller: S. Perry

Description: Soil Boring

Logged By: S. Hubbs

Contractor: New England Geotech Borehole Dia.: 2.00in

Date(s): 11/30/06 - 11/30/06

Ground Elevation: 0.00' Coordinate X: 0.00

Drilling Method: Geoprobe

Total Depth: 8.00'

Coordinate Y: 0.00

Back Fill:

type: Native Material

fm: 0.00'

to: 8.00'

Remarks: Field Instrument: Photovac 2020 Refusal at 8.0 feet.

type:

fm:

to:

fm:

				type: fm: type: fm: type: fm:		to: to: to:	
Elevation	Depth	Sample No.	Recovery	Material Description	Graphic Log	USCS Code	РНОТ
-	_	-11 N/A -12		0-0.5': SAND, F; some gravel; dusky brown (5YR 2/2), dry. 0.5-1.1': SAND, M; coal and ash; dusky brown (5YR 2/2), dry. 1.1-5.0': Sand, M-C and gravel; dusky brown (5YR 2/2), moist.		SW FI	O ppm
-2-	2-	N/A			0.000		
-4-	4	 N/A			0.0	SW	0 ppm
-6-	6 —			Sand, M and gravel; moderate brown (5YR 4/4), wet.	0 0		
-8-	8-			End of boring at 8.0 feet.	0		
-10-	10 -						
-12-	12-						
-	-						
-14 -	14 —						
-16 -	16-						
-18-	18- ed By:			P	age 1	of 1	

Total Depth: 7.50'

Project Location: Franklin, Massachusetts

Remarks: Field Instrument: Photovac 2020

Refusal at 7.5 feet.

Site Id: B-10

Project Number: 2005-0458 B10



Location: See map Datum: Logged By: S. Hubbs Description: Soil Boring Ground Elevation: 0.00' Date(s): 12/01/06 - 12/01/06 Coordinate X: 0.00

Driller: S. Perry Contractor: New England Geotech Borehole Dia.: 2.00in

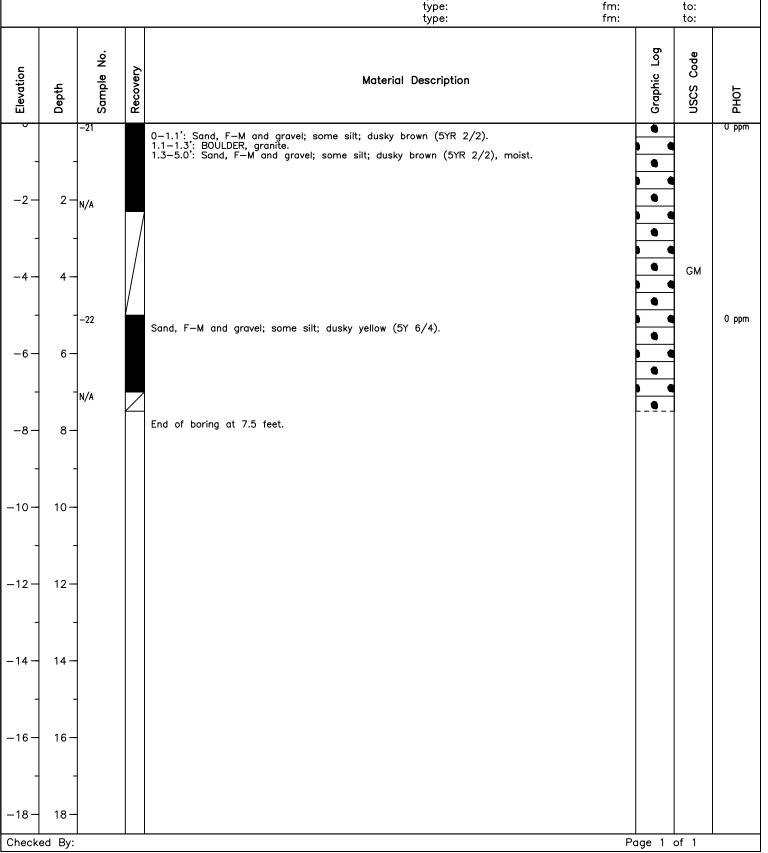
Drilling Method: Geoprobe

Back Fill:

Coordinate Y: 0.00 type: Native Material

type: type: fm: 0.00' to: 7.50' fm: to:

fm: to: fm: to: fm: to:



Project Location: Franklin, Massachusetts

Refusal at 8.0 feet.

Site Id: B-11

Project Number: 2005-0458 B10



Location: See map Description: Soil Boring Datum:

Logged By: S. Hubbs Driller: S. Perry Contractor: New England Geotech Borehole Dia.: 2.00in

Ground Elevation: 0.00' Date(s): 12/01/06 - 12/01/06 Coordinate X: 0.00 Drilling Method: Geoprobe

Total Depth: 8.00'

Coordinate Y: 0.00

Remarks: Field Instrument: Photovac 2020

Back Fill:

type: Native Material fm: 0.00'

to: 8.00' fm: to: fm: to:

type:

type:

				type: 1m: type: fm: type: fm:		to: to: to:	
Elevation	Depth	Sample No.	Recovery	Material Description	Graphic Log	USCS Code	РНОТ
-	-	-23		0-1.6': SAND, F-M, yellowish gray (5Y 7/2) to dusky brown (5YR 2/2) at 0.4 feet, dry. 1.6-2.4': SAND, F; some silt; trace gravel; yellowish gray (5Y 7/2), moist. 2.4-2.7': Sand, F and silt; wood or textile; moderate brown (5YR 3/4), moist. 2.7-5.0': SAND, F; some silt; trace gravel; yellowish gray (5Y 7/2), moist.	0.00	SW	O ppm
-2-	2 -	N/A					
-4-	4 —						
-6-	6 —	-24	,	Sand, F and silt; dusky brown (5YR 3/4), moist to wet at 6.2 feet.		SM	0 ppm
-8-	- 8 <i>-</i>	N/A					
-	-			End of boring at 8.0 feet.			
-10 <i>-</i>	10 -						
-12-	12-						
-14-	14 —						
-16-	16-						
	-						
-18 -	18 – ed Bv:			· · · · · · · · · · · · · · · · · · ·	Page 1	of 1	
LONGER	Ju Dy.			<u>'</u>	29C	~' '	

Project Location: Franklin, Massachusetts

Site Id: B-12

Project Number: 2005-0458 B10



Location: See map

Datum:

Logged By: S. Hubbs

Driller: S. Perry

to: 8.00'

to:

Description: Soil Boring

Total Depth: 8.00'

Ground Elevation: 0.00'

Contractor: New England Geotech Borehole Dia.: 2.00in

Date(s): 12/01/06 - 12/01/06

Coordinate X: 0.00 Coordinate Y: 0.00

Back Fill:

Remarks: Field Instrument: Photovac 2020

Refusal at 8.0 feet.

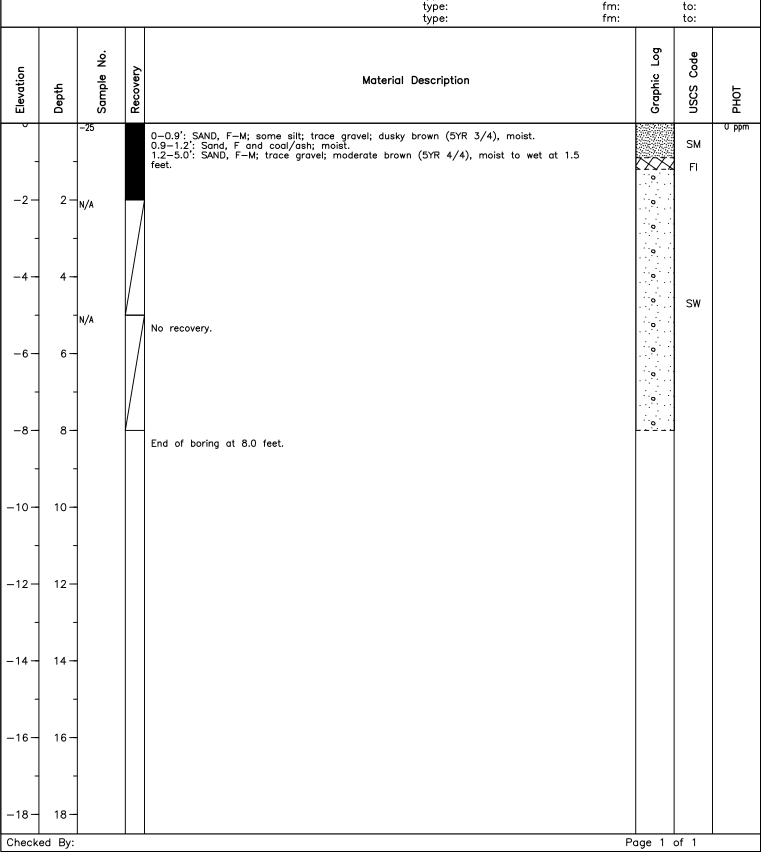
type: Native Material

Drilling Method: Geoprobe

fm: 0.00'

type: fm: type: fm:

to: fm: to: fm: to:





APPENDIX B

FORMER NU-STYLE COMPANY, INC. PHASE II SITE ASSESSMENT

PREMIER LABORATORY CERTIFICATES OF ANALYSIS, FUSS & O'NEILL DATA VERIFICATION NARRATIVES AND CERTIFICATIONS, AND DATA VALIDATION COMPLETION WORKSHEETS



Modified Tier I Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

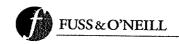
Premier Laboratory Project Number:	E612024
Date Samples Received at Laboratory:	12/1/2006
Date of Review:	1/11/2007

Seven soil samples were collected and submitted to Premier Laboratory, LLC in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant metals plus barium by EPA Methods 6010B and 7471, cyanide by EPA Method 9012, polychlorinated biphenyls (PCBs) by EPA Method 8082, and petroleum hydrocarbons by Massachusetts Department of Environmental Protection (MADEP) Methods Extractable Petroleum Hydrocarbons (EPH) and Volatile Petroleum Hydrocarbons (VPH). One aqueous trip blank was also submitted for analysis of VOCs by EPA Method 8260B. Dedicated sampling equipment was employed; therefore, no equipment blank was indicated. A field duplicate was collected and submitted during the first day of sampling.

Samples were analyzed within method-specified holding times and in accordance with the Massachusetts Contingency Plan (MCP) Compendium of Analytical Methods (CAM) data enhancement protocols.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

	<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREMENTS:	/	
Field measurement calibration records	□⁄	
Groundwater field measurements (if applicable)		
Soil sampling field measurements (if applicable)		
Sediment sampling field measurements (if applicable)		
Surface water sampling field measurements (if applicable)		□ N/A
Low-flow sampling field measurements (if applicable)	Π,	A4 [
Documentation of field activities	Ø,	
Sample numbering and labeling		
Chain-of-Custody records	☑,	
Trip blanks		
Duplicate samples		□ NA
Equipment blanks		DNA
Split samples (if any)		o NA
2. LABORATORY MEASUREMENTS:		
Trip blanks	Ø	
Instrument blanks	<u> </u>	
Laboratory control samples	$\overline{\square}$	
Duplicates samples		NA
Equipment blanks		- NA
Matrix spike/matrix spike duplicates		□ NIA
Analysis type		
Chain-of-Custody records	_ ☑_	
Surrogate recoveries	<u> </u>	
Sample Project Narratives		П
Split samples (if any)		□ NA
TOTAL: _	13	0



Premier Laboratory, LLC 61 Louisa Viens Drive Dayville, CT 06241 Telephone: 860-774-6814 Fax: 860-774-26A9

ANALYTICAL DATA & QUALITY CONTROL **REPORT**

Report Number: E612024

Project: 20050458.B10/Nu-Style Phase II

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss





ANALYTICAL DATA REPORT

Report Number: E612024 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

Received Date: 12/1/2006 Report Date: 3/13/2007

Premier Laboratory, LLC Authorized Signature



956B850

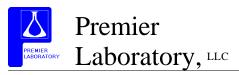
Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

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MADEP MCP Analytical Method Report Certification Form										
Labo	ratory Name: Prer	nier Laboratory, LL	С		Project #:	E61202	24			
Proje	ct Location: Frank	klin, MA			MADEP RT	N ¹ :				
	This Form provides certifications for the following data set:[list Laboratory Sample ID Number(s)] 1, 2, 3, 4, 5, 6, 7, 8									
Samp	Sample Matrices: ☐ Groundwater ☒ Soil/Sediment ☐ Drinking Water ☐ Other									
MCP	SW-846	8260B ⊠	8151A □	8330	□ 6010	OB ⊠		7470	0A/1A ⊠	
Meth	ods Used	8270C □	8081A □	VPH		20 🗆		90	14M ² □	
	ecified in MADEP	8082 ⊠	8021B □	EPH	⊠ 7000 \$	S ³ □		7	7196A 🗆	
Analyt	endium of ical Methods. all that apply)	2 M - SW-846 Method	ng Number (RTN), if know 9014 or MADEP Physiol s 7000 Series List individ	ogically Available		Method				
	An affirmative res	ponse to question	ns A, B, C, and D is	required for	"Presumptive	Certair	ity"	sta	tus	
А	•	•	ooratory in a condition dy documentation fo			⊠ Ye	es		No ¹	
В	Were all QA/QC p	rocedures required	I for the specified an	alytical metho	d(s)	⊠ Ye	es		No ¹	
	included in this rep	oort followed, includ	ding the requirement	to note and						
	discuss in a narrat	tive QC data that di	id not meet appropria	ate performan	ce					
	standards or guide	elines?								
С	Does the analytica	al data included in t	this report meet all th	ne requiremer	nts	⊠ Ye	es		No ¹	
	for "Presumptive C	Certainty", as descr	ibed in Section 2.0 (a),(b),(c) and	(d) of the					
	MADEP documen	t CAM VII A, "Qual	ity Assurance and C	uality Control	Guidelines					
	•	and Reporting of A								
D			the VPH or EPH me	thod run with	out	⊠ Ye	es		No ¹	
	significant modific	ations, as specified	d in Section 11.3?							
	_	•	nd F below is requir		umptive Certa					
E	Were all QC perfo specified methods		and recommendatior	ns for the		⊠ Ye	3 S		No ¹	
F	Were results for a	Il analyte-list comp	ounds/elements for t	he specified		⊠ Ye	es		No ¹	
	method(s) reported	d?								
	¹ All NO answ	ers must be addres	ssed in an attached	Environmenta	l Laboratory ca	ase narra	ative	Э.		
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.										
Sign	ature:	*Elekkin		Position:	Laboratory Di	rector				
Print	ed Name: Rober	t Stevenson		Date: <u>3/13</u>	3/2007					

Page 2 of 46



> Report No: E612024 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory received eight samples from Fuss & O'Neill on 12/01/2006. The samples were analyzed from the following list of analyses:

Cyanide, Total, by 9012 in GW/SW 9012[9012]
Mercury by 7471 in SW 7471[7471]
Trace Priority Pollutant (13) Metals in Soil 6010B[3000], 7471[7471]
Volatiles by 8260B in GW/SW 8260B

Extractable Petroleum Hydrocarbon (EPH) MADEP EPH[MADEP EPH] Moisture, Percent PCB's by 8082 in GW/SW 8082[3500] Volatile Petroleum Hydrocarbon (VPH) MADEP VPH

In order to meet requested detection limits, EDB results were estimated to 3 ppb for EPA method 8260B. Dibromochloromethane, 1,2-Dichlorobenzene and 1,1,2,2-Tetrachloroethane were all estimated to a value of 5.0 ppb. This value of 5.0 ppb corresponds to the lowest level of calibration on the instrument prior to the % solid value being calculated into the reported detection limits. The samples were ND for all estimated compounds to their respective values.

Variances:

SDG:

A full list 8260B LCS was run and met the applicable recovery criteria for "Presumptive Certainty". An LCS Duplicate encompasing all target compounds was not run for EPA method 8260B. Both an LCS and LCSD were analyzed for the oxygenate compounds only.

Method:

None reported.

QA/QC:

956B850

Sample 1A, 841061201-19, Volatiles by 8260B: Three internal standards were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 1A, 841061201-19, Volatiles by 8260B: Two surrogate spikes were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

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> Report No: E612024 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY (continued)

QA/QC (continued):

Sample 2A, 841061201-20, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 5A, 841061201-23, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 5A, 841061201-23, Volatiles by 8260B: One surrogate spike was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

Sample 6A, 841061201-24, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 7A, 841061201-25, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 7A, 841061201-25, Volatiles by 8260B: Two surrogate spikes were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

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Laboratory: Premier Laboratory, LLC
PL Report No: E612024
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(1) 841061201-19					
Date Collected: 12/1/2006 Matrix	· Solid				
Cyanide, Total, by SW-846 9012	ND	0.54	mg/kg	12/06/06 1	2:02 DDD
Trace Metals by 6010B	112	0.5 1		12,00,00	2.02
Antimony	6.5	0.50	mg/kg	12/05/06	BSZ
Arsenic	3.1	0.50	mg/kg	12/05/06	BSZ
Barium	110	0.50	mg/kg	12/05/06	BSZ
Beryllium	0.37	0.050	mg/kg	12/05/06	BSZ
Cadmium	0.54	0.10	mg/kg	12/05/06	BSZ
Chromium	27	0.50	mg/kg	12/05/06	BSZ
Copper	29	0.50	mg/kg	12/05/06	BSZ
Lead	780	2.2	mg/kg	12/05/06	BSZ 10
Nickel	6.4	0.50	mg/kg	12/05/06	BSZ
Selenium	ND	0.50	mg/kg	12/05/06	BSZ
Silver	ND	0.10	mg/kg	12/05/06	BSZ
Thallium	ND	0.25	mg/kg	12/05/06	BSZ
Zinc	310	0.50	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	0.073	0.022	mg/kg	12/05/06	AM
(2) 841061201-20					
Date Collected: 12/1/2006 Matrix	: Solid				
Cyanide, Total, by SW-846 9012	ND	0.58	mg/kg	12/06/06 1	2:03 DDD
Trace Metals by 6010B					
Antimony	6.9	0.50	mg/kg	12/05/06	BSZ
Arsenic	ND	0.50	mg/kg	12/05/06	BSZ
Barium	55	0.50	mg/kg	12/05/06	BSZ
Beryllium	0.17	0.050	mg/kg	12/05/06	BSZ
Cadmium	0.18	0.10	mg/kg	12/05/06	BSZ
Chromium	26	0.50	mg/kg	12/05/06	BSZ
Copper	9.5	0.50	mg/kg	12/05/06	BSZ
Lead	310	0.20	mg/kg	12/05/06	BSZ
Nickel	6.3	0.50	mg/kg	12/05/06	BSZ
Selenium	ND	0.50	mg/kg	12/05/06	BSZ
Silver	ND	0.10	mg/kg	12/05/06	BSZ
Thallium	ND	0.25	mg/kg	12/05/06	BSZ
Zinc	84	0.50	mg/kg	12/05/06	BSZ

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Laboratory: Premier Laboratory, LLC

PL Report No: E612024

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(2) 841061201-20 (continued) Date Collected: 12/1/2006 M Mercury by SW-846 7471 in SW	<u>(atrix: Solid</u> ND	0.023	mg/kg	12/05/06	AM
(3) 841061201-21					
	<u> Iatrix: Solid</u>				
Cyanide, Total, by SW-846 9012	ND	0.55	mg/kg	12/06/06 12:04	DDD
Trace Metals by 6010B					
Antimony	ND	0.50	mg/kg	12/05/06	BSZ
Arsenic	ND	0.50	mg/kg	12/05/06	BSZ
Barium	16	0.50	mg/kg	12/05/06	BSZ
Beryllium	0.70	0.050	mg/kg	12/05/06	BSZ
Cadmium	ND	0.10	mg/kg	12/05/06	BSZ
Chromium	5.2	0.50	mg/kg	12/05/06	BSZ
Copper	6.3	0.50	mg/kg	12/05/06	BSZ
Lead	2.9	0.20	mg/kg	12/05/06	BSZ
Nickel	3.6	0.50	mg/kg	12/05/06	BSZ
Selenium	ND	0.50	mg/kg	12/05/06	BSZ
Silver	ND	0.10	mg/kg	12/05/06	BSZ
Thallium	ND	0.27	mg/kg	12/05/06	BSZ
Zinc	22	0.50	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	0.023	0.022	mg/kg	12/05/06	AM
(4) 841061201-22 <u>Date Collected: 12/1/2006</u> M Cyanide, Total, by SW-846 9012	<u>(atrix: Solid</u> ND	0.54	mg/kg	12/06/06 12:07	DDD

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Laboratory: Premier Laboratory, LLC
PL Report No: E612024
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(4) 841061201-22 (continued)					
Date Collected: 12/1/2006 Matrix: Soli	d				
Trace Metals by 6010B	<u>u</u>				
Antimony	ND	0.50	mg/kg	12/05/06	BSZ
Anumony Arsenic	ND	0.50	mg/kg	12/05/06	BSZ
Barium	9.6	0.50	mg/kg	12/05/06	BSZ
Beryllium	0.91	0.050	mg/kg	12/05/06	BSZ
Cadmium	ND	0.030	mg/kg	12/05/06	BSZ
Chromium	1.9	0.10	mg/kg	12/05/06	BSZ
Copper	1.9	0.50	mg/kg	12/05/06	BSZ
Lead	4.8	0.30	mg/kg	12/05/06	BSZ
Nickel	1.0	0.20	mg/kg	12/05/06	BSZ
Selenium	ND	0.50	mg/kg	12/05/06	BSZ
Silver	ND ND	0.30	mg/kg	12/05/06	BSZ
					BSZ
Thallium Zinc	ND	0.27 0.50	mg/kg	12/08/06	BSZ BSZ
	15 ND		mg/kg	12/05/06	
Mercury by SW-846 7471 in SW	ND	0.021	mg/kg	12/05/06	AM
(5) 841061201-23					
Date Collected: 12/1/2006 Matrix: Soli	a.				
	_	0.54	ma a /lra	12/06/06 12	2:08 DDD
Cyanide, Total, by SW-846 9012	ND	0.54	mg/kg	12/06/06 12	2:08 DDD
Trace Metals by 6010B	ND	0.54	ma a /lra	12/05/06	BSZ
Antimony	ND	0.54	mg/kg	12/05/06	BSZ BSZ
Arsenic	1.8	0.54	mg/kg	12/05/06	
Barium	23	0.54	mg/kg	12/05/06	BSZ
Beryllium	0.16	0.054	mg/kg	12/05/06	BSZ
Cadmium	0.26	0.11	mg/kg	12/08/06	BSZ
Chromium	5.4	0.54	mg/kg	12/05/06	BSZ
Copper	8.5	0.54	mg/kg	12/05/06	BSZ
Lead	17	0.22	mg/kg	12/05/06	BSZ
Nickel	3.2	0.54	mg/kg	12/05/06	BSZ
Selenium	ND	0.54	mg/kg	12/05/06	BSZ
Silver	ND	0.11	mg/kg	12/05/06	BSZ
Thallium	ND	0.27	mg/kg	12/05/06	BSZ
Zinc	48	0.54	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	0.032	0.022	mg/kg	12/05/06	AM

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Laboratory: Premier Laboratory, LLC
PL Report No: E612024
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter		Result	DL	Units	Completed		By	Dilution
(6) 841061201-24								
Date Collected: 12/1/2006	Matrix: Solid							
Cyanide, Total, by SW-846 9012	Matria: Bond	ND	0.6	mg/kg	12/06/06	12:09	DDD)
Trace Metals by 6010B		T\D	0.0	mg/kg	12/00/00	12.07	DDL	
Antimony		ND	0.60	mg/kg	12/05/06		BSZ	
Arsenic		ND	0.60	mg/kg	12/05/06		BSZ	
Barium		34	0.60	mg/kg	12/05/06		BSZ	
Beryllium		0.38	0.060	mg/kg	12/05/06		BSZ	
Cadmium		ND	0.12	mg/kg	12/05/06		BSZ	
Chromium		4.4	0.60	mg/kg	12/05/06		BSZ	
Copper		2.9	0.60	mg/kg	12/05/06		BSZ	
Lead		4.3	0.24	mg/kg	12/05/06		BSZ	
Nickel		1.7	0.60	mg/kg	12/05/06		BSZ	
Selenium		ND	0.60	mg/kg	12/05/06		BSZ	
Silver		ND	0.12	mg/kg	12/05/06		BSZ	
Thallium		ND	0.30	mg/kg	12/08/06		BSZ	
Zinc		8.4	0.60	mg/kg	12/05/06		BSZ	
Mercury by SW-846 7471 in SW		ND	0.024	mg/kg	12/05/06		AM	
(7) 841061201-25								
Date Collected: 12/1/2006	Matrix: Solid							
Cyanide, Total, by SW-846 9012		ND	0.56	mg/kg	12/06/06	12:10	DDD)
Trace Metals by 6010B								
Antimony		ND	0.56	mg/kg	12/08/06		BSZ	
Arsenic		3.1	0.56	mg/kg	12/05/06		BSZ	
Barium		30	0.56	mg/kg	12/05/06		BSZ	
Beryllium		0.26	0.056	mg/kg	12/05/06		BSZ	
Cadmium		0.19	0.11	mg/kg	12/05/06		BSZ	
Chromium		6.0	0.56	mg/kg	12/05/06		BSZ	
Copper		37	0.56	mg/kg	12/05/06		BSZ	
Lead		93	0.22	mg/kg	12/05/06		BSZ	
Nickel		130	0.56	mg/kg	12/05/06		BSZ	
Selenium		ND	0.56	mg/kg	12/05/06		BSZ	
Silver		ND	0.11	mg/kg	12/05/06		BSZ	
Thallium		ND	0.28	mg/kg	12/05/06		BSZ	
Zinc		28	0.56	mg/kg	12/05/06		BSZ	

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Laboratory: Premier Laboratory, LLC

PL Report No: E612024

Date Received: 12/1/2006

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II

 Parameter
 Result
 DL
 Units
 Completed
 By Dilution

 (7) 841061201-25 (continued)
 Date Collected: 12/1/2006

 Mercury by SW-846 7471 in SW
 0.044
 0.022
 mg/kg
 12/05/06
 AM

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061201-19
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/04/06	Percent Moisture:	8.2
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6400	ug/kg
C9-C12 Aliphatics**	50	14000	6400	ug/kg
C9-C10 Aromatics***	50	ND	6400	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	96	70%-130%
2,5-dibromotoluene #2	104	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	320	ug/kg
Ethylbenzene	ND	320	ug/kg
Methyl tert-butyl ether (MTBE)	ND	64	ug/kg
Naphthalene	ND	320	ug/kg
Toluene	ND	320	ug/kg
m,p-Xylenes	ND	320	ug/kg
o-Xylene	ND	320	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

VOLATILE ORGANIC ANALYSIS DATA SHEET

Matrix: Solid

Percent Moisture: 8.2

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 Sample Description: 841061201-19

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By: GPDilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28525.D;J28740.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	18
71-43-2	Benzene	ND	4.6
108-86-1	Bromobenzene	ND	4.6
74-97-5	Bromochloromethane	ND	4.6
75-27-4	Bromodichloromethane	ND	4.6
75-25-2	Bromoform	ND	4.6
74-83-9	Bromomethane	ND	9.2
78-93-3	2-Butanone (MEK)	ND	9.2
104-51-8	n-Butylbenzene	ND	4.6
135-98-8	sec-Butylbenzene	ND	4.6
98-06-6	tert-Butylbenzene	ND	4.6
75-15-0	Carbon disulfide	ND	4.6
56-23-5	Carbon tetrachloride	ND	4.6
108-90-7	Chlorobenzene	ND	4.6
75-00-3	Chloroethane	ND	9.2
67-66-3	Chloroform	ND	4.6
74-87-3	Chloromethane	ND	9.2
95-49-8	2-Chlorotoluene	ND	4.6
106-43-4	4-Chlorotoluene	ND	4.6
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	4.6
124-48-1	Dibromochloromethane	ND	4.6
106-93-4	1,2-Dibromoethane (EDB)	ND	2.8
74-95-3	Dibromomethane	ND	4.6
95-50-1	1,2-Dichlorobenzene	ND	4.6
541-73-1	1,3-Dichlorobenzene	ND	4.6
106-46-7	1,4-Dichlorobenzene	ND	4.6
75-71-8	Dichlorodifluoromethane	ND	9.2
75-34-3	1,1-Dichloroethane	ND	4.6
107-06-2	1,2-Dichloroethane	ND	4.6
75-35-4	1,1-Dichloroethene	ND	4.6
156-59-2	cis-1,2-Dichloroethene	ND	4.6
156-60-5	trans-1,2-Dichloroethene	ND	4.6
78-87-5	1,2-Dichloropropane	ND	4.6
142-28-9	1,3-Dichloropropane	ND	4.6
590-20-7	2,2-Dichloropropane	ND	4.6
563-58-6	1,1-Dichloropropene	ND	4.6
10061-01-5	cis-1,3-Dichloropropene	ND	4.6
10061-02-6	trans-1,3-Dichloropropene	ND	4.6
60-29-7	Diethyl ether	ND	9.2
123-91-1	1,4-Dioxane	ND	18

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VOLATILE ORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 (continued) Sample Description: 841061201-19

Date Collected:12/1/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:8.2Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28525.D;J28740.D

Units: ug/kg

CAS No.	Parameter	F	Result	DL
	Ethyl tertiary-butyl ether (EtBE)		ND	50
100-41-4	Ethylbenzene		ND	4.6
87-68-3	Hexachlorobutadiene		ND	4.6
591-78-6	2-Hexanone		ND	9.2
98-82-8	Isopropylbenzene		ND	4.6
99-87-6	4-Isopropyltoluene		ND	4.6
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	4.6
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	9.2
75-09-2	Methylene chloride		ND	4.6
91-20-3	Naphthalene		ND	4.6
103-65-1	n-Propylbenzene		ND	4.6
100-42-5	Styrene		ND	4.6
994-05-8	Tertiary-amyl methyl ether (TAME)		ND	50
109-99-9	Tetrahydrofuran		ND	4.6
96-18-4	1,2,3-Trichloropropane		ND	4.6
630-20-6	1,1,1,2-Tetrachloroethane		ND	4.6
79-34-5	1,1,2,2-Tetrachloroethane		ND	4.6
127-18-4	Tetrachloroethene (PCE)		ND	4.6
108-88-3	Toluene		ND	4.6
87-61-6	1,2,3-Trichlorobenzene		ND	4.6
120-82-1	1,2,4-Trichlorobenzene		ND	4.6
71-55-6	1,1,1-Trichloroethane		ND	4.6
79-00-5	1,1,2-Trichloroethane		ND	4.6
79-01-6	Trichloroethene (TCE)		ND	4.6
75-69-4	Trichlorofluoromethane		ND	9.2
95-63-6	1,2,4-Trimethylbenzene		ND	4.6
108-67-8	1,3,5-Trimethylbenzene		ND	4.6
75-01-4	Vinyl chloride		ND	9.2
95-47-6	o-Xylene		ND	4.6
	m,p-Xylenes		ND	4.6
Surrogate	Recovery	Limits		
Bromofluorobenzene	61%	78%-111%		
1,2-Dichloroethane-d4	108%	91%-114%		
Toluene-d8	118%	86%-115%		

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061201-20
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/04/06	Percent Moisture:	13.7
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7600	ug/kg
C9-C12 Aliphatics**	50	ND	7600	ug/kg
C9-C10 Aromatics***	50	ND	7600	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	89	70%-130%
2,5-dibromotoluene #2	98	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	380	ug/kg
Ethylbenzene	ND	380	ug/kg
Methyl tert-butyl ether (MTBE)	ND	76	ug/kg
Naphthalene	ND	380	ug/kg
Toluene	ND	380	ug/kg
m,p-Xylenes	ND	380	ug/kg
o-Xylene	ND	380	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

VOLATILE ORGANIC ANALYSIS DATA SHEET

Matrix: Solid

Percent Moisture: 13.7 Sample Weight/Volume:

Dilution Factor: 1

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Legation: Frontlin MA

Location: Franklin, MA
PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 2 Sample Description: 841061201-20

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/04/06 By: GP

Method: 8260B Soil Extract Volume:
QC Batch#: 50061 Lab Data File: J28501.D;J28741.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	22
71-43-2	Benzene	ND	5.6
108-86-1	Bromobenzene	ND	5.6
74-97-5	Bromochloromethane	ND	5.6
75-27-4	Bromodichloromethane	ND	5.6
75-25-2	Bromoform	ND	5.6
74-83-9	Bromomethane	ND	11
78-93-3	2-Butanone (MEK)	ND	11
104-51-8	n-Butylbenzene	ND	5.6
135-98-8	sec-Butylbenzene	ND	5.6
98-06-6	tert-Butylbenzene	ND	5.6
75-15-0	Carbon disulfide	ND	5.6
56-23-5	Carbon tetrachloride	ND	5.6
108-90-7	Chlorobenzene	ND	5.6
75-00-3	Chloroethane	ND	11
67-66-3	Chloroform	ND	5.6
74-87-3	Chloromethane	ND	11
95-49-8	2-Chlorotoluene	ND	5.6
106-43-4	4-Chlorotoluene	ND	5.6
108-20-3	Di-isopropyl ether (DIPE)	ND	56
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.6
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.4
74-95-3	Dibromomethane	ND	5.6
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.6
106-46-7	1,4-Dichlorobenzene	ND	5.6
75-71-8	Dichlorodifluoromethane	ND	11
75-34-3	1,1-Dichloroethane	ND	5.6
107-06-2	1,2-Dichloroethane	ND	5.6
75-35-4	1,1-Dichloroethene	ND	5.6
156-59-2	cis-1,2-Dichloroethene	ND	5.6
156-60-5	trans-1,2-Dichloroethene	ND	5.6
78-87-5	1,2-Dichloropropane	ND	5.6
142-28-9	1,3-Dichloropropane	ND	5.6
590-20-7	2,2-Dichloropropane	ND	5.6
563-58-6	1,1-Dichloropropene	ND	5.6
10061-01-5	cis-1,3-Dichloropropene	ND	5.6
10061-02-6	trans-1,3-Dichloropropene	ND	5.6
60-29-7	Diethyl ether	ND	11
123-91-1	1,4-Dioxane	ND	22

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VOLATILE ORGANIC ANALYSIS DATA SHEET

Matrix: Solid

Percent Moisture: 13.7 Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 (continued) Sample Description: 841061201-20

Date Collected: 12/1/2006
Date Received: 12/1/2006
Date Extracted: By:
Date Analyzed: 12/04/06 By: GP

Date Analyzed: 12/04/06 By: GP Dilution Factor: 1
Method: 8260B Soil Extract Volume:

Units: ug/kg

QC Batch#: 50061 Lab Data File: J28501.D;J28741.D

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	56
100-41-4	Ethylbenzene	ND	5.6
87-68-3	Hexachlorobutadiene	ND	5.6
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.6
99-87-6	4-Isopropyltoluene	ND	5.6
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.6
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.6
91-20-3	Naphthalene	ND	5.6
103-65-1	n-Propylbenzene	ND	5.6
100-42-5	Styrene	ND	5.6
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	56
109-99-9	Tetrahydrofuran	ND	5.6
96-18-4	1,2,3-Trichloropropane	ND	5.6
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.6
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	ND	5.6
108-88-3	Toluene	ND	5.6
87-61-6	1,2,3-Trichlorobenzene	ND	5.6
120-82-1	1,2,4-Trichlorobenzene	ND	5.6
71-55-6	1,1,1-Trichloroethane	ND	5.6
79-00-5	1,1,2-Trichloroethane	ND	5.6
79-01-6	Trichloroethene (TCE)	ND	5.6
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.6
108-67-8	1,3,5-Trimethylbenzene	ND	5.6
75-01-4	Vinyl chloride ND		11
95-47-6	· · · · · · · · · · · · · · · · · · ·		5.6
	m,p-Xylenes	ND	5.6
Surrogate	Recovery Limits		

Surrogate	Recovery	Limits
Bromofluorobenzene	85%	78%-111%
1,2-Dichloroethane-d4	101%	91%-114%
Toluene-d8	105%	86%-115%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061201-21
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/04/06	Percent Moisture:	8.5
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6400	ug/kg
C9-C12 Aliphatics**	50	ND	6400	ug/kg
C9-C10 Aromatics***	50	ND	6400	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	88	70%-130%
2,5-dibromotoluene #2	98	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	320	ug/kg
Ethylbenzene	ND	320	ug/kg
Methyl tert-butyl ether (MTBE)	ND	64	ug/kg
Naphthalene	ND	320	ug/kg
Toluene	ND	320	ug/kg
m,p-Xylenes	ND	320	ug/kg
o-Xylene	ND	320	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes *** Excludes Naphthalene

VOLATILE ORGANIC ANALYSIS DATA SHEET

Matrix: Solid

Percent Moisture: 8.5 Sample Weight/Volume:

Dilution Factor: 50

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill

Laboratory: Franklin MA

Location: Franklin, MA
PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 3 Sample Description: 841061201-21

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Analyzed: 12/05/06 By: GP

Method: 8260B Soil Extract Volume:
QC Batch#: 50078 Soil Extract Volume:
Lab Data File: M32494.D;J28742.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	1100
71-43-2	Benzene	ND	270
108-86-1	Bromobenzene	ND	270
74-97-5	Bromochloromethane	ND	270
75-27-4	Bromodichloromethane	ND	270
75-25-2	Bromoform	ND	270
74-83-9	Bromomethane	ND	550
78-93-3	2-Butanone (MEK)	ND	550
104-51-8	n-Butylbenzene	ND	270
135-98-8	sec-Butylbenzene	ND	270
98-06-6	tert-Butylbenzene	ND	270
75-15-0	Carbon disulfide	ND	270
56-23-5	Carbon tetrachloride	ND	270
108-90-7	Chlorobenzene	ND	270
75-00-3	Chloroethane	ND	550
67-66-3	Chloroform	ND	270
74-87-3	Chloromethane	ND	550
95-49-8	2-Chlorotoluene	ND	270
106-43-4	4-Chlorotoluene	ND	270
108-20-3	Di-isopropyl ether (DIPE)	ND	55
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	270
124-48-1	Dibromochloromethane	ND	270
106-93-4	1,2-Dibromoethane (EDB)	ND	160
74-95-3	Dibromomethane	ND	270
95-50-1	1,2-Dichlorobenzene	ND	270
541-73-1	1,3-Dichlorobenzene	ND	270
106-46-7	1,4-Dichlorobenzene	ND	270
75-71-8	Dichlorodifluoromethane	ND	550
75-34-3	1,1-Dichloroethane	ND	270
107-06-2	1,2-Dichloroethane	ND	270
75-35-4	1,1-Dichloroethene	ND	270
156-59-2	cis-1,2-Dichloroethene	ND	270
156-60-5	trans-1,2-Dichloroethene	ND	270
78-87-5	1,2-Dichloropropane	ND	270
142-28-9	1,3-Dichloropropane	ND	270
590-20-7	2,2-Dichloropropane	ND	270
563-58-6	1,1-Dichloropropene	ND	270
10061-01-5	cis-1,3-Dichloropropene	ND	270
10061-02-6	trans-1,3-Dichloropropene	ND	270
60-29-7	Diethyl ether	ND	550
123-91-1	1,4-Dioxane	ND	1100

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VOLATILE ORGANIC ANALYSIS DATA SHEET

Matrix: Solid

Percent Moisture: 8.5 Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 (continued) Sample Description: 841061201-21

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/05/06 By: GP

Date Analyzed: 12/05/06 By: GP Dilution Factor: 50
Method: 8260B Soil Extract Volume:
QC Batch#: 50078 Lab Data File: M32494.D;J28742.D

Units: ug/kg

CAS No.	Parameter		DL
	Ethyl tertiary-butyl ether (EtBE)	ND	55
100-41-4	Ethylbenzene	ND	270
87-68-3	Hexachlorobutadiene	ND	270
591-78-6	2-Hexanone	ND	550
98-82-8	Isopropylbenzene	ND	270
99-87-6	4-Isopropyltoluene	ND	270
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	270
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	550
75-09-2	Methylene chloride	ND	270
91-20-3	Naphthalene	ND	270
103-65-1	n-Propylbenzene	ND	270
100-42-5	Styrene	ND	270
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	55
109-99-9	Tetrahydrofuran	ND	270
96-18-4	1,2,3-Trichloropropane	ND	270
630-20-6	1,1,1,2-Tetrachloroethane	ND	270
79-34-5	1,1,2,2-Tetrachloroethane	ND	270
127-18-4	Tetrachloroethene (PCE)	4300	270
108-88-3	Toluene	ND	270
87-61-6	1,2,3-Trichlorobenzene	ND	270
120-82-1	1,2,4-Trichlorobenzene	ND	270
71-55-6	1,1,1-Trichloroethane	ND	270
79-00-5	1,1,2-Trichloroethane	ND	270
79-01-6	Trichloroethene (TCE)	9300	270
75-69-4	Trichlorofluoromethane	ND	550
95-63-6	1,2,4-Trimethylbenzene	ND	270
108-67-8	1,3,5-Trimethylbenzene	ND	270
75-01-4	Vinyl chloride	ND	550
95-47-6	o-Xylene	ND	270
	m,p-Xylenes	ND	270
Surrogate	Recovery	Limits	
Bromofluorobenzene	88%	78%-111%	
1,2-Dichloroethane-d4	102%	91%-114%	

Toluene-d8 105% 86%-115%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061201-22
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/04/06	Percent Moisture:	6.6
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6100	ug/kg
C9-C12 Aliphatics**	50	ND	6100	ug/kg
C9-C10 Aromatics***	50	ND	6100	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	91	70%-130%
2,5-dibromotoluene #2	101	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	300	ug/kg
Ethylbenzene	ND	300	ug/kg
Methyl tert-butyl ether (MTBE)	ND	61	ug/kg
Naphthalene	ND	300	ug/kg
Toluene	ND	300	ug/kg
m,p-Xylenes	ND	300	ug/kg
o-Xylene	ND	300	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

VOLATILE ORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 Sample Description: 841061201-22

Date Collected: 12/1/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 6.6

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50075 Lab Data File: J28527.D;J28743.D

Units: ug/kg

Acetone	CAS No.	Parameter	Result	DL
108-86-1 Bromobenzene	67-64-1	Acetone	ND	22
74-97-5 Bromochloromethane ND 5.4 75-27-4 Bromodichloromethane ND 5.4 75-25-2 Bromoform ND 5.4 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.4 135-98-8 sec-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 80-6-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chloroform ND 5.4 74-87-3 Chloroform ND 5.4 74-87-3 Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3	71-43-2	Benzene	ND	5.4
75-27-4 Bromoform ND 5.4 75-25-2 Bromoform ND 5.4 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 18-93-3 1-Butylbenzene ND 5.4 135-98-8 sec-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 108-90-7 Chlorobenzene ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorobenzene ND 5.4 78-03-3 Chloroform ND 5.4 78-79-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 109-34	108-86-1	Bromobenzene	ND	5.4
75-25-2 Bromoform ND 5.4 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.4 135-98-8 sec-Butylbenzene ND 5.4 135-98-8 sec-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 56-23-5 Carbon disulfide ND 5.4 56-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorotofum ND 5.4 75-00-3 Chlorotofum ND 5.4 74-87-3 Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 106-34-1 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 106-93-4 1,2-Dibromo-dhoromethane ND 5.0	74-97-5	Bromochloromethane	ND	5.4
74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 54 135-98-8 sec-Butylbenzene ND 54 98-06-6 tert-Butylbenzene ND 54 75-15-0 Carbon disulfide ND 54 56-23-5 Carbon tetrachloride ND 54 108-90-7 Chlorobenzene ND 54 75-00-3 Chloroftenane ND 11 67-66-3 Chloroform ND 54 74-87-3 Chlorotoluene ND 54 106-43-4 4-Chlorotoluene ND 54 108-20-3 Di-isopropyl ether (DIPE) ND 54 108-20-3 Di-isopropyl ether (DIPE) ND 54 124-48-1 Dibromochloromethane ND 50 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 54 14-95-3 Dibromochloromethane ND 50	75-27-4	Bromodichloromethane	ND	5.4
78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.4 135-98-8 sec-Burylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 56-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorochane ND 11 67-66-3 Chloroform ND 5.4 74-87-3 Chloromethane ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 104-34-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 106-93-4 1,2-Dichoromethane ND 5.0 12-48-1 Dibromochlane (EDB) ND 5.4 14-95-3 Dibromochlane (EDB) ND 5.4 <	75-25-2	Bromoform	ND	5.4
78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.4 135-98-8 sec-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 56-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chloroform ND 5.4 75-00-3 Chloroform ND 5.4 74-87-3 Chloroform ND 5.4 106-43-4 Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 108-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 74-95-3 Dibromochloromethane ND 5.0 106-46-7 1,2-Dichlorobenzene ND	74-83-9	Bromomethane	ND	11
135-98-8 sec-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 56-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorothane ND 11 67-66-3 Chloroform ND 5.4 74-87-3 Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 104-93-1 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 105-93-2 1,2-Dibromo-stane (EDB) ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene	78-93-3		ND	11
135-98-8 sec-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 108-90-7 Chlorobenzene ND 5.4 108-90-7 Chlorobenzene ND 5.4 108-90-7 Chlorothane ND 11 67-66-3 Chlorothane ND 11 95-49-8 Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 106-93-4 1,2-Dibromoethane ND 5.0 106-93-4 1,2-Dibromoethane EDB ND 3.2 149-5-3 Dibromoethane EDB ND 5.4 195-50-1 1,2-Dichlorobenzene ND 5.4 108-46-7 1,4-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 107-06-5 trans-1,2-Dichlorothene ND 5.4 156-50-5 trans-1,2-Dichlorothene ND 5.4 156-60-5 trans-1,2-Dichlorothene ND 5.4 142-28-9 1,3-Dichlorothene ND 5.4 156-35-8-6 1,1-Dichloropopane ND 5.4 10061-01-5 cis-1,3-Dichloropopene ND 5.4 10061-02-6 trans-1,3-Dichloropopene ND 5.4 10061-02-6	104-51-8	n-Butylbenzene	ND	5.4
98-06-6 tert-Butylbenzene ND 5.4 75-15-0 Carbon disulfide ND 5.4 75-15-0 Carbon disulfide ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorotehane ND 11 67-66-3 Chloroform ND 5.4 74-87-3 Chlorotehane ND 5.4 96-48-8 2-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 106-28 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 106-93-4 1,2-Dibromo-chloromethane ND 5.0 14-48-1 Dibromo-chloromethane (EDB) ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene	135-98-8		ND	5.4
75-15-0 Carbon disulfide ND 5.4 56-23-5 Carbon tetrachloride ND 5.4 108-90-7 Chlorobenzene ND 5.4 75-00-3 Chlorochtane ND 11 67-66-3 Chloroform ND 5.4 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 108-21-4 1.2-Dichlorothonomethane ND 5.0 109-18-4 1.2-Dichlorothonomethane <td>98-06-6</td> <td>· · · · · · · · · · · · · · · · · · ·</td> <td>ND</td> <td>5.4</td>	98-06-6	· · · · · · · · · · · · · · · · · · ·	ND	5.4
108-90-7	75-15-0	•	ND	5.4
75-00-3 Chloroethane ND 11 67-66-3 Chloroform ND 5.4 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochlane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 107-46-7 1,4-Dichlorobenzene ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroptone ND	56-23-5	Carbon tetrachloride	ND	5.4
67-66-3 Chloroform ND 5.4 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 54 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 107-06-2 1,2-Dichlorothene ND 5.4 156-60-5 trans-1,2-Dichlorothene ND 5.4 156-60-5 trans-1,2-Dichloropropane	108-90-7	Chlorobenzene	ND	5.4
74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 5.4 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,2-Dichlorotethane ND 5.4 107-06-2 1,2-Dichlorotethane ND 5.4 156-69-2 cis-1,2-Dichlorotethene ND 5.4 156-60-5 trans-1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropro	75-00-3	Chloroethane	ND	11
95-49-8 2-Chlorotoluene ND 5.4 106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 54 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloropropane ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 500-20-7 2,2-Dichloro	67-66-3	Chloroform	ND	5.4
106-43-4 4-Chlorotoluene ND 5.4 108-20-3 Di-isopropyl ether (DIPE) ND 54 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromoethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 5063-58-6 1,1-Di	74-87-3	Chloromethane	ND	11
108-20-3 Di-isopropyl ether (DIPE) ND 54 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 15-71-8 Dichlorodifluoromethane ND 5.4 15-34-3 1,1-Dichlorothane ND 5.4 107-06-2 1,2-Dichlorothane ND 5.4 156-59-2 cis-1,2-Dichlorothene ND 5.4 156-60-5 trans-1,2-Dichlorothene ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 506-38-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,	95-49-8	2-Chlorotoluene	ND	5.4
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloropropane ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-02-6 trans-	106-43-4	4-Chlorotoluene	ND	5.4
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.4 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethene ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroptopene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 503-58-6 1,1-Dichlo	108-20-3	Di-isopropyl ether (DIPE)	ND	54
106-93-4 1,2-Dibromoethane (EDB) ND 3.2 74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether	96-12-8		ND	5.4
74-95-3 Dibromomethane ND 5.4 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 5.4	124-48-1	Dibromochloromethane	ND	5.0
95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
541-73-1 1,3-Dichlorobenzene ND 5.4 106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	74-95-3	Dibromomethane	ND	5.4
106-46-7 1,4-Dichlorobenzene ND 5.4 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	95-50-1	1,2-Dichlorobenzene	ND	5.0
75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	541-73-1	1,3-Dichlorobenzene	ND	5.4
75-34-3 1,1-Dichloroethane ND 5.4 107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	106-46-7	1,4-Dichlorobenzene	ND	5.4
107-06-2 1,2-Dichloroethane ND 5.4 75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	75-71-8	Dichlorodifluoromethane	ND	11
75-35-4 1,1-Dichloroethene ND 5.4 156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	75-34-3	1,1-Dichloroethane	ND	5.4
156-59-2 cis-1,2-Dichloroethene ND 5.4 156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	107-06-2	1,2-Dichloroethane	ND	5.4
156-60-5 trans-1,2-Dichloroethene ND 5.4 78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	75-35-4	1,1-Dichloroethene	ND	5.4
78-87-5 1,2-Dichloropropane ND 5.4 142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	156-59-2	cis-1,2-Dichloroethene	ND	5.4
142-28-9 1,3-Dichloropropane ND 5.4 590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	156-60-5	trans-1,2-Dichloroethene	ND	5.4
590-20-7 2,2-Dichloropropane ND 5.4 563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	78-87-5	1,2-Dichloropropane	ND	5.4
563-58-6 1,1-Dichloropropene ND 5.4 10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	142-28-9	1,3-Dichloropropane	ND	5.4
10061-01-5 cis-1,3-Dichloropropene ND 5.4 10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	590-20-7	2,2-Dichloropropane	ND	5.4
10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	563-58-6		ND	5.4
10061-02-6 trans-1,3-Dichloropropene ND 5.4 60-29-7 Diethyl ether ND 11	10061-01-5	cis-1,3-Dichloropropene	ND	5.4
60-29-7 Diethyl ether ND 11	10061-02-6	trans-1,3-Dichloropropene	ND	5.4
123-91-1 1,4-Dioxane ND 22	60-29-7		ND	11
	123-91-1	1,4-Dioxane	ND	22

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612024 PL Sample No: 4 (continued) Sample Description: 841061201-22

Date Collected: 12/1/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 6.6 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/05/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume:

Lab Data File: J28527.D;J28743.D QC Batch#: 50075

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	54
100-41-4	Ethylbenzene	ND	5.4
87-68-3	Hexachlorobutadiene	ND	5.4
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.4
99-87-6	4-Isopropyltoluene	ND	5.4
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.4
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.4
91-20-3	Naphthalene	ND	5.4
103-65-1	n-Propylbenzene	ND	5.4
100-42-5	Styrene	ND	5.4
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	54
109-99-9	Tetrahydrofuran	ND	5.4
96-18-4	1,2,3-Trichloropropane	ND	5.4
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.4
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	48	5.4
108-88-3	Toluene	ND	5.4
87-61-6	1,2,3-Trichlorobenzene	ND	5.4
120-82-1	1,2,4-Trichlorobenzene	ND	5.4
71-55-6	1,1,1-Trichloroethane	ND	5.4
79-00-5	1,1,2-Trichloroethane	ND	5.4
79-01-6	Trichloroethene (TCE)	150	5.4
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.4
108-67-8	1,3,5-Trimethylbenzene	ND	5.4
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.4
	m,p-Xylenes	ND	5.4
Surrogate	Recovery	Limits	
Bromofluorobenzene	94%	78%-111%	
1,2-Dichloroethane-d4	102%	91%-114%	
TP 1 10	0.507	0.60/ 1150/	

Surrogate	Recovery	Limits
Bromofluorobenzene	94%	78%-111%
1,2-Dichloroethane-d4	102%	91%-114%
Toluene-d8	95%	86%-115%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061201-23
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	7.8
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	12000	ug/kg
C9-C12 Aliphatics**	50	ND	12000	ug/kg
C9-C10 Aromatics***	50	ND	12000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	91	70%-130%
2,5-dibromotoluene #2	101	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	620	ug/kg
Ethylbenzene	ND	620	ug/kg
Methyl tert-butyl ether (MTBE)	ND	120	ug/kg
Naphthalene	ND	620	ug/kg
Toluene	ND	620	ug/kg
m,p-Xylenes	ND	620	ug/kg
o-Xylene	ND	620	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841061201-23

Date Collected:12/1/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:7.8Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28528.D;J28744.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	20
71-43-2	Benzene	ND	5.0
108-86-1	Bromobenzene	ND	5.0
74-97-5	Bromochloromethane	ND	5.0
75-27-4	Bromodichloromethane	ND	5.0
75-25-2	Bromoform	ND	5.0
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.0
135-98-8	sec-Butylbenzene	ND	5.0
98-06-6	tert-Butylbenzene	ND	5.0
75-15-0	Carbon disulfide	ND	5.0
56-23-5	Carbon tetrachloride	ND	5.0
108-90-7	Chlorobenzene	ND	5.0
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.0
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.0
106-43-4	4-Chlorotoluene	ND	5.0
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.0
74-95-3	Dibromomethane	ND	5.0
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.0
106-46-7	1,4-Dichlorobenzene	ND	5.0
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.0
107-06-2	1,2-Dichloroethane	ND	5.0
75-35-4	1,1-Dichloroethene	ND	5.0
156-59-2	cis-1,2-Dichloroethene	ND	5.0
156-60-5	trans-1,2-Dichloroethene	ND	5.0
78-87-5	1,2-Dichloropropane	ND	5.0
142-28-9	1,3-Dichloropropane	ND	5.0
590-20-7	2,2-Dichloropropane	ND	5.0
563-58-6	1,1-Dichloropropene	ND	5.0
10061-01-5	cis-1,3-Dichloropropene	ND	5.0
10061-02-6	trans-1,3-Dichloropropene	ND	5.0
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	20

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Matrix: Solid

Percent Moisture: 7.8 Sample Weight/Volume:

Dilution Factor: 1

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 (continued) Sample Description: 841061201-23

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/05/06 By: GP

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28528.D;J28744.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	50
100-41-4	Ethylbenzene	ND	5.0
87-68-3	Hexachlorobutadiene	ND	5.0
591-78-6	2-Hexanone	ND	10
98-82-8	Isopropylbenzene	ND	5.0
99-87-6	4-Isopropyltoluene	ND	5.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	5.0
103-65-1	n-Propylbenzene	ND	5.0
100-42-5	Styrene	ND	5.0
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	50
109-99-9	Tetrahydrofuran	ND	5.0
96-18-4	1,2,3-Trichloropropane	ND	5.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	40	5.0
108-88-3	Toluene	16	5.0
87-61-6	1,2,3-Trichlorobenzene	ND	5.0
120-82-1	1,2,4-Trichlorobenzene	ND	5.0
71-55-6	1,1,1-Trichloroethane	ND	5.0
79-00-5	1,1,2-Trichloroethane	ND	5.0
79-01-6	Trichloroethene (TCE)	5.0	5.0
75-69-4	Trichlorofluoromethane	ND	10
95-63-6	1,2,4-Trimethylbenzene	ND	5.0
108-67-8	1,3,5-Trimethylbenzene	ND	5.0
75-01-4	Vinyl chloride	ND	10
95-47-6	o-Xylene	ND	5.0
	m,p-Xylenes	ND	5.0
Surrogate	Recovery Lir	nits	
Bromofluorobenzene	72% 78°	%-111%	

Surrogate	Recovery	Limits
Bromofluorobenzene	72%	78%-111%
1,2-Dichloroethane-d4	104%	91%-114%
Toluene-d8	115%	86%-115%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061201-24
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	16.4
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	8000	ug/kg
C9-C12 Aliphatics**	50	ND	8000	ug/kg
C9-C10 Aromatics***	50	ND	8000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	95	70%-130%
2,5-dibromotoluene #2	104	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	400	ug/kg
Ethylbenzene	ND	400	ug/kg
Methyl tert-butyl ether (MTBE)	ND	80	ug/kg
Naphthalene	ND	400	ug/kg
Toluene	ND	400	ug/kg
m,p-Xylenes	ND	400	ug/kg
o-Xylene	ND	400	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Custome

PL Report No: E612024 PL Sample No: 6

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Analyzed: 12/05/06 By: GP

Method: 8260B QC Batch#: 50075

Units: ug/kg

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061201-24

Matrix: Solid

Percent Moisture: 16.4 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume:

Lab Data File: J28529.D;J28745.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	30	23
71-43-2	Benzene	ND	5.7
108-86-1	Bromobenzene	ND	5.7
74-97-5	Bromochloromethane	ND	5.7
75-27-4	Bromodichloromethane	ND	5.7
75-25-2	Bromoform	ND	5.7
74-83-9	Bromomethane	ND	11
78-93-3	2-Butanone (MEK)	ND	11
104-51-8	n-Butylbenzene	ND	5.7
135-98-8	sec-Butylbenzene	ND	5.7
98-06-6	tert-Butylbenzene	ND	5.7
75-15-0	Carbon disulfide	ND	5.7
56-23-5	Carbon tetrachloride	ND	5.7
108-90-7	Chlorobenzene	ND	5.7
75-00-3	Chloroethane	ND	11
67-66-3	Chloroform	ND	5.7
74-87-3	Chloromethane	ND	11
95-49-8	2-Chlorotoluene	ND	5.7
106-43-4	4-Chlorotoluene	ND	5.7
108-20-3	Di-isopropyl ether (DIPE)	ND	57
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.7
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.4
74-95-3	Dibromomethane	ND	5.7
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.7
106-46-7	1,4-Dichlorobenzene	ND	5.7
75-71-8	Dichlorodifluoromethane	ND	11
75-34-3	1,1-Dichloroethane	ND	5.7
107-06-2	1,2-Dichloroethane	ND	5.7
75-35-4	1,1-Dichloroethene	ND	5.7
156-59-2	cis-1,2-Dichloroethene	ND	5.7
156-60-5	trans-1,2-Dichloroethene	ND	5.7
78-87-5	1,2-Dichloropropane	ND	5.7
142-28-9	1,3-Dichloropropane	ND	5.7
590-20-7	2,2-Dichloropropane	ND	5.7
563-58-6	1,1-Dichloropropene	ND	5.7
10061-01-5	cis-1,3-Dichloropropene	ND	5.7
10061-02-6	trans-1,3-Dichloropropene	ND	5.7
60-29-7	Diethyl ether	ND	11
123-91-1	1,4-Dioxane	ND	23

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Matrix: Solid

Percent Moisture: 16.4

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 6 (continued) Sample Description: 841061201-24

Date Collected: 12/1/2006
Date Received: 12/1/2006
Date Extracted: By:

Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By: GPDilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28529.D;J28745.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	57
100-41-4	Ethylbenzene	ND	5.7
87-68-3	Hexachlorobutadiene	ND	5.7
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.7
99-87-6	4-Isopropyltoluene	ND	5.7
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.7
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.7
91-20-3	Naphthalene	ND	5.7
103-65-1	n-Propylbenzene	ND	5.7
100-42-5	Styrene	ND	5.7
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	57
109-99-9	Tetrahydrofuran	ND	5.7
96-18-4	1,2,3-Trichloropropane	ND	5.7
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.7
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	45	5.7
108-88-3	Toluene	ND	5.7
87-61-6	1,2,3-Trichlorobenzene	ND	5.7
120-82-1	1,2,4-Trichlorobenzene	ND	5.7
71-55-6	1,1,1-Trichloroethane	ND	5.7
79-00-5	1,1,2-Trichloroethane	ND	5.7
79-01-6	Trichloroethene (TCE)	ND	5.7
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.7
108-67-8	1,3,5-Trimethylbenzene	ND	5.7
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.7
	m,p-Xylenes	ND	5.7
Surrogate	Recovery Lir	mits	
Bromofluorobenzene	78% 78'	%-111%	-

Surrogate	Recovery	Limits
Bromofluorobenzene	78%	78%-111%
1,2-Dichloroethane-d4	102%	91%-114%
Toluene-d8	105%	86%-115%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	7	Sample Description:	841061201-25
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	10.2
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6800	ug/kg
C9-C12 Aliphatics**	50	ND	6800	ug/kg
C9-C10 Aromatics***	50	ND	6800	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	102	70%-130%
2,5-dibromotoluene #2	112	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	340	ug/kg
Ethylbenzene	ND	340	ug/kg
Methyl tert-butyl ether (MTBE)	ND	68	ug/kg
Naphthalene	ND	340	ug/kg
Toluene	ND	340	ug/kg
m,p-Xylenes	ND	340	ug/kg
o-Xylene	ND	340	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill

Laboratory: Franklin MA

Location: Franklin, MA
PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 7 Sample Description: 841061201-25

Date Collected: 12/1/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Analyzed: 12/05/06 By: GP

Method: 8260B QC Batch#: 50075

Units: ug/kg

Matrix: Solid Percent Moisture: 10.2 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume:

Lab Data File: J28530.D;J28746.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	21
71-43-2	Benzene	ND	5.2
108-86-1	Bromobenzene	ND	5.2
74-97-5	Bromochloromethane	ND	5.2
75-27-4	Bromodichloromethane	ND	5.2
75-25-2	Bromoform	ND	5.2
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.2
135-98-8	sec-Butylbenzene	ND	5.2
98-06-6	tert-Butylbenzene	ND	5.2
75-15-0	Carbon disulfide	ND	5.2
56-23-5	Carbon tetrachloride	ND	5.2
108-90-7	Chlorobenzene	ND	5.2
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.2
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.2
106-43-4	4-Chlorotoluene	ND	5.2
108-20-3	Di-isopropyl ether (DIPE)	ND	52
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.2
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.1
74-95-3	Dibromomethane	ND	5.2
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.2
106-46-7	1,4-Dichlorobenzene	ND	5.2
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.2
107-06-2	1,2-Dichloroethane	ND	5.2
75-35-4	1,1-Dichloroethene	ND	5.2
156-59-2	cis-1,2-Dichloroethene	ND	5.2
156-60-5	trans-1,2-Dichloroethene	ND	5.2
78-87-5	1,2-Dichloropropane	ND	5.2
142-28-9	1,3-Dichloropropane	ND	5.2
590-20-7	2,2-Dichloropropane	ND	5.2
563-58-6	1,1-Dichloropropene	ND	5.2
10061-01-5	cis-1,3-Dichloropropene	ND	5.2
10061-02-6	trans-1,3-Dichloropropene	ND	5.2
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	21

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Matrix: Solid

Percent Moisture: 10.2 Sample Weight/Volume:

Dilution Factor: 1

Customer: Fuss & O'Neill Laboratory: Premier Laboratory, LLC Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 (continued) Sample Description: 841061201-25

Date Collected: 12/1/2006 Date Received: 12/1/2006 Date Extracted: By: Date Analyzed: 12/05/06 By: GP

Method: 8260B Soil Extract Volume: Lab Data File: J28530.D;J28746.D QC Batch#: 50075

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	52
100-41-4	Ethylbenzene	ND	5.2
87-68-3	Hexachlorobutadiene	ND	5.2
591-78-6	2-Hexanone	ND	10
98-82-8	Isopropylbenzene	ND	5.2
99-87-6	4-Isopropyltoluene	ND	5.2
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.2
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10
75-09-2	Methylene chloride	ND	5.2
91-20-3	Naphthalene	ND	5.2
103-65-1	n-Propylbenzene	ND	5.2
100-42-5	Styrene	ND	5.2
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	52
109-99-9	Tetrahydrofuran	ND	5.2
96-18-4	1,2,3-Trichloropropane	ND	5.2
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.2
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	11	5.2
108-88-3	Toluene	ND	5.2
87-61-6	1,2,3-Trichlorobenzene	ND	5.2
120-82-1	1,2,4-Trichlorobenzene	ND	5.2
71-55-6	1,1,1-Trichloroethane	ND	5.2
79-00-5	1,1,2-Trichloroethane	ND	5.2
79-01-6	Trichloroethene (TCE)	6.5	5.2
75-69-4	Trichlorofluoromethane	ND	10
95-63-6	1,2,4-Trimethylbenzene	ND	5.2
108-67-8	1,3,5-Trimethylbenzene	ND	5.2
75-01-4	Vinyl chloride	ND	10
95-47-6	o-Xylene	ND	5.2
	m,p-Xylenes	ND	5.2
Surrogate	Recovery	Limits	
Bromofluorobenzene	68%	78%-111%	
1,2-Dichloroethane-d4	96%	91%-114%	

Toluene-d8 121% 86%-115%

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 Sample Description: 841061201-26

Date Collected: 12/1/2006 Matrix: Solid

Date Received:12/1/2006Percent Moisture:N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By: GPDilution Factor:50Method:8260BSoil Extract Volume:

QC Batch#: 51012 Lab Data File: M32507.D;M32616.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	250
71-43-2	Benzene	ND	50
108-86-1	Bromobenzene	ND	50
74-97-5	Bromochloromethane	ND	50
75-27-4	Bromodichloromethane	ND	50
75-25-2	Bromoform	ND	50
74-83-9	Bromomethane	ND	50
78-93-3	2-Butanone (MEK)	ND	250
104-51-8	n-Butylbenzene	ND	50
135-98-8	sec-Butylbenzene	ND	50
98-06-6	tert-Butylbenzene	ND	50
75-15-0	Carbon disulfide	ND	50
56-23-5	Carbon tetrachloride	ND	50
108-90-7	Chlorobenzene	ND	50
75-00-3	Chloroethane	ND	50
67-66-3	Chloroform	ND	50
74-87-3	Chloromethane	ND	50
95-49-8	2-Chlorotoluene	ND	50
106-43-4	4-Chlorotoluene	ND	50
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	50
124-48-1	Dibromochloromethane	ND	50
106-93-4	1,2-Dibromoethane (EDB)	ND	50
74-95-3	Dibromomethane	ND	50
95-50-1	1,2-Dichlorobenzene	ND	50
541-73-1	1,3-Dichlorobenzene	ND	50
106-46-7	1,4-Dichlorobenzene	ND	50
75-71-8	Dichlorodifluoromethane	ND	50
75-34-3	1,1-Dichloroethane	ND	50
107-06-2	1,2-Dichloroethane	ND	50
75-35-4	1,1-Dichloroethene	ND	50
156-59-2	cis-1,2-Dichloroethene	ND	50
156-60-5	trans-1,2-Dichloroethene	ND	50
78-87-5	1,2-Dichloropropane	ND	50
142-28-9	1,3-Dichloropropane	ND	50
590-20-7	2,2-Dichloropropane	ND	50
563-58-6	1,1-Dichloropropene	ND	50
10061-01-5	cis-1,3-Dichloropropene	ND	50
10061-02-6	trans-1,3-Dichloropropene	ND	50
60-29-7	Diethyl ether	ND	50
123-91-1	1,4-Dioxane	ND	1000

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 (continued) Sample Description: 841061201-26

Date Collected: 12/1/2006 Matrix: Solid

Date Received:12/1/2006Percent Moisture:N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By: GPDilution Factor:50Method:8260BSoil Extract Volume:

QC Batch#: 51012 Lab Data File: M32507.D;M32616.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	50
100-41-4	Ethylbenzene	ND	50
87-68-3	Hexachlorobutadiene	ND	50
591-78-6	2-Hexanone	ND	250
98-82-8	Isopropylbenzene	ND	50
99-87-6	4-Isopropyltoluene	ND	50
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	50
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	250
75-09-2	Methylene chloride	ND	50
91-20-3	Naphthalene	ND	50
103-65-1	n-Propylbenzene	ND	50
100-42-5	Styrene	ND	50
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	50
109-99-9	Tetrahydrofuran	ND	50
96-18-4	1,2,3-Trichloropropane	ND	50
630-20-6	1,1,1,2-Tetrachloroethane	ND	50
79-34-5	1,1,2,2-Tetrachloroethane	ND	50
127-18-4	Tetrachloroethene (PCE)	ND	50
108-88-3	Toluene	ND	50
87-61-6	1,2,3-Trichlorobenzene	ND	50
120-82-1	1,2,4-Trichlorobenzene	ND	50
71-55-6	1,1,1-Trichloroethane	ND	50
79-00-5	1,1,2-Trichloroethane	ND	50
79-01-6	Trichloroethene (TCE)	ND	50
75-69-4	Trichlorofluoromethane	ND	50
95-63-6	1,2,4-Trimethylbenzene	ND	50
108-67-8	1,3,5-Trimethylbenzene	ND	50
75-01-4	Vinyl chloride	ND	50
95-47-6	o-Xylene	ND	50
	m,p-Xylenes	ND	50
Surrogate	Recovery Lir	nits	
Bromofluorobenzene	93% 879	%-105%	

Surrogate	Recovery	Limits
Bromofluorobenzene	93%	87%-105%
1,2-Dichloroethane-d4	94%	91%-109%
Toluene-d8	104%	92%-105%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061201-19
Preservative	None	-	
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	8.2
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	52	40%-140%
2-Bromonaphthalene	68	40%-140%
2-Fluorobiphenyl	69	40%-140%
o-Terphenyl	40	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 1 Sample Description: 841061201-19

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 8.2

Date Extracted: 12/07/06 By: AKB Sample Weight/Volume: 30.07 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#:

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120822.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	67%	30%-150%
Decachlorobiphenyl	74%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061201-20
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	13.7
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	53	40%-140%
2-Bromonaphthalene	77	40%-140%
2-Fluorobiphenyl	80	40%-140%
o-Terphenyl	52	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 Sample Description: 841061201-20

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moist

Date Received: 12/1/2006 Percent Moisture: 13.7

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.01 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#:

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120836.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15
C .	T)	т • •,	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	53%	30%-150%
Decachlorobiphenyl	55%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061201-21
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	8.5
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	ND	10000	ug/kg
C11-C22 Aromatics*	1	25000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	78	40%-140%
2-Bromonaphthalene	63	40%-140%
2-Fluorobiphenyl	66	40%-140%
o-Terphenyl	58	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	200	100	ug/kg
Anthracene	250	100	ug/kg
Benzo[a]anthracene	1500	100	ug/kg
Benzo[a]pyrene	2000	100	ug/kg
Benzo[b]fluoranthene	2000	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	860	100	ug/kg
Chrysene	120	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	1300	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	260	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	260	100	ug/kg
Pyrene	1600	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 Sample Description: 841061201-21

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 8.5

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.31 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#:

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120837.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14
Surrogate	Pacovary	Limite	_

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	50%	30%-150%
Decachlorobiphenyl	53%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061201-22
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	6.6
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	ND	10000	ug/kg
C11-C22 Aromatics*	1	ND	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	55	40%-140%
2-Bromonaphthalene	66	40%-140%
2-Fluorobiphenyl	70	40%-140%
o-Terphenyl	40	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	ND	100	ug/kg
Anthracene	ND	100	ug/kg
Benzo[a]anthracene	ND	100	ug/kg
Benzo[a]pyrene	ND	100	ug/kg
Benzo[b]fluoranthene	ND	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	ND	100	ug/kg
Chrysene	ND	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	ND	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	ND	100	ug/kg
Pyrene	ND	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 Sample Description: 841061201-22

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moist

Units: ug/kg

Date Received: 12/1/2006 Percent Moisture: 6.6
Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.11 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120840.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	59%	30%-150%
Decachlorobiphenyl	53%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061201-23
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/11/06	Percent Moisture:	7.8
Date Analyzed:	12/12/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	18000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	59	40%-140%
2-Bromonaphthalene	65	40%-140%
2-Fluorobiphenyl	69	40%-140%
o-Terphenyl	49	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841061201-23

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 7.8

Units: ug/kg

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.08 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120841.D

 CAS No.
 Parameter
 Result
 DL

 12674-11-2
 Aroclor 1016
 ND
 14

 11104-28-2
 Aroclor 1221
 ND
 14

 11141-16-5
 Aroclor 1232
 ND
 14

11141-16-5 Aroclor 1242 53469-21-9 ND 14 Aroclor 1248 14 12672-29-6 ND 11097-69-1 Aroclor 1254 ND 14 Aroclor 1260 14 11096-82-5 ND

SurrogateRecoveryLimitsTetrachloro-m-xylene84%30%-150%Decachlorobiphenyl114%30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061201-24
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	16.4
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	69	40%-140%
2-Bromonaphthalene	62	40%-140%
2-Fluorobiphenyl	64	40%-140%
o-Terphenyl	60	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 6 Sample Description: 841061201-24

Aroclor 1260

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moist

Units: ug/kg

11096-82-5

Date Received: 12/1/2006 Percent Moisture: 16.4

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.78 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120842.D

CAS No. Parameter Result DL Aroclor 1016 12674-11-2 ND 16 11104-28-2 Aroclor 1221 ND 16 11141-16-5 Aroclor 1232 ND 16 Aroclor 1242 53469-21-9 ND 16 Aroclor 1248 12672-29-6 ND 16 11097-69-1 Aroclor 1254 ND 16

SurrogateRecoveryLimitsTetrachloro-m-xylene72%30%-150%Decachlorobiphenyl53%30%-150%

ND

16

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612024	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	7	Sample Description:	841061201-25
Preservative	None		
		Dilution (Target):	1
Date Collected:	12/1/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/01/06	Percent Moisture:	10.2
Date Analyzed:	12/07/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	16000	11000	ug/kg
C11-C22 Aromatics*	1	28000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	45	40%-140%
2-Bromonaphthalene	71	40%-140%
2-Fluorobiphenyl	72	40%-140%
o-Terphenyl	42	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	120	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	490	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	290	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	1300	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	940	110	ug/kg
Pyrene	1200	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612024 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 Sample Description: 841061201-25

Date Collected: 12/1/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 10.2

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.23 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120843.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	65%	30%-150%
Decachlorobiphenyl	61%	30%-150%

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Thin H. Wyros 22°C

3

□ 146 Hartford Road, Manchester, CT 06040

☐ 56 Quarry Road, Trumbull, CT 06611

□ 1419 Richland Street, Columbia, SC 29201

□ 78 Interstate Drive, West Springfield, MA 01089

□ 610 Lynndale Court, Suite E, Greenville, NC 27858 □ 24 Madison Avenue Extension, Albany, NY 12203

Additional Comments:

- See attache QA/QC checkist

275 Promenade Street, Suite 350, Providence, RI 02908

□ 80 Washington Street, Suite 301, Poughkeepsie, NY 12001

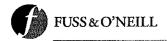
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Modified Tier I Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E612529
Date Samples Received at Laboratory:	12/11/2006
Date of Review:	1/11/2007

Six aqueous samples, including one field duplicate, were collected by low-flow methodology and submitted to Premier Laboratory, LLC in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant metals plus barium by EPA Methods 6010B and 7471, and petroleum hydrocarbons by Massachusetts Department of Environmental Protection (MADEP) Methods Extractable Petroleum Hydrocarbons (EPH) and Volatile Petroleum Hydrocarbons (VPH). One aqueous trip blank was also submitted for analysis of VOCs by EPA Method 8260B. Dedicated sampling equipment was employed; therefore, no equipment blank was indicated.

Samples were analyzed within method-specified holding times and in accordance with the Massachusetts Contingency Plan (MCP) Compendium of Analytical Methods (CAM) data enhancement protocols.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

4 CAMPI INC AND PIELD MEASUREMENT	YES	NO NO
1. SAMPLING AND FIELD MEASUREMEN		_
Field measurement calibration records	<u> </u>	/ 🗆
Groundwater field measurements (if applicable)		
Soil sampling field measurements (if applicable)		□ 'Y/
Sediment sampling field measurements (if applicable)		
Surface water sampling field measurements (if applicable	e)	
Low-flow sampling field measurements (if applicable)		, 🗆
Documentation of field activities		
Sample numbering and labeling	Image: section of the content of the	
Chain-of-Custody records	9	
Trip blanks		
Duplicate samples		
Equipment blanks		□ NA
Split samples (if any)		
2. LABORATORY MEASUREMENTS:	,	
Trip blanks	V	
Instrument blanks		
Laboratory control samples		
Duplicates samples	$\overline{\mathbf{Z}}$	
Equipment blanks	$\overline{\Box}$	TNA
Matrix spike/matrix spike duplicates	$\overline{\Box}$	_ ~ / <i>A</i>
Analysis type	ল	n
Chain-of-Custody records		n
Surrogate recoveries	ra/	
Sample Project Narratives	الم	ä
Split samples (if any)		. HAIA
opin samples (if any)		□ <i>[/</i> 4
TO	OTAL:	
PE	ERCENT COMPLETE:	%



Premier Laboratory, LLC 61 Louisa Viens Drive Dayville, CT 06241

Telephone: 860-774-6814 Fax: 860-774-26A9

ANALYTICAL DATA & QUALITY CONTROL REPORT

Report Number: E612529

Project: 20050458.B10/Nu-Style Phase II

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss





61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

ANALYTICAL DATA REPORT

Report Number: E612529 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

Received Date: 12/11/2006 Report Date: 12/19/2006

Premier Laboratory, LLC Authorized Signature

7. Wank



958B150

Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

Page 1 of 32



61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

MADEP MCP Analytical Method Report Certification Form									
Laboratory Name: Premier Laboratory, LLC Project #:						E61252	29		
Proje	Project Location: Franklin, MA MADEP RTN ¹ :								
	•	fications for the foll	lowing data set:[list Li	aboratory Sar	mple ID Numbe	er(s)]			
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			2.7/0.7/2						
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	SW-846		8151A 🗆	8330 VDU		0B ⊠			A/1A ⊠ 14M ² □
	ods Used	8270C □ 8082 □	8081A □ 8021B □	VPH EPH		20 🗆			196A □
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Analyt	ical Methods.	2 M - SW-846 Method	19014 or MADEP Physiol	ogically Available		Method			
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С			ribed in Section 2.0 (a	•			:5	ш 1	NO
	•		lity Assurance and Q		` '				
		and Reporting of	-	tuality Contro	Guidelines				
D			the VPH or EPH me	athod run with	out	⊠ Ye			No ¹
		ations, as specified		tilloa rair witii	out	Б 10	.3	ш,	10
			nd F below is requir	red for "Pres	umntive Certa	inty" sta	atus		
E	•	•	and recommendation		difficate derivation	∏ Ye			Vo ¹
	specified methods		and recommendation	10 101 1110		、	,,	_ :	
F	· •		ounds/elements for t	the specified		⊠ Ye	es		No ¹
	method(s) reported			·					
			ssed in an attached i	Environmenta	al Laboratory ca	ase narra	ative),	
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this									
			nd belief, accurate and c						
	-								
Sign	ature:	M. Which		Position:	Laboratory Dir	rector			
Print	ted Name: Ronald	d Warila		Date: <u>12/</u>	19/2006				

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61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

> Report No: E612529 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory received seven samples from Fuss & O'Neill on 12/11/2006. The samples were analyzed from the following list of analyses:

Extractable Petroleum Hydrocarbon (EPH)
MADEP EPH[MADEP EPH]
Volatile Petroleum Hydrocarbon (VPH)
MADEP VPH

Trace Priority Pollutant (13) Metals in Water 6010B[3000], 7470A[245.1] Volatiles by 8260B (GA/GW-1/S-1) 8260B

Variances:

SDG:

None reported.

Method:

None reported.

QA/QC:

Sample 1C, 841061208-27, Volatiles by 8260B (GA/GW-1/S-1): One surrogate spike was outside quality control limits for the matrix spike/matrix spike duplicate due to matrix interference. All surrogate recoveries were within limits for the sample.

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INORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, LLC
PL Report No: E612529
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/11/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(1) 941071309 37					
(1) 841061208-27 Date Collected: 12/8/2006 Matrix: Aqueous					
<u>-</u>					
Trace Metals by 6010B	ND	0.010	ma/I	12/18/06	BSZ
Antimony Arsenic	ND ND	0.010	mg/L	12/18/06	BSZ
Barium	0.042	0.010	mg/L mg/L	12/18/06	BSZ
Beryllium	0.042 ND	0.0010	mg/L mg/L	12/18/06	BSZ
Cadmium	ND ND	0.0010	mg/L mg/L	12/18/06	BSZ
Chromium	ND ND	0.0020	mg/L mg/L	12/18/06	BSZ
Copper	ND ND	0.010	mg/L mg/L	12/18/06	BSZ
Lead	0.014	0.010	_	12/18/06	BSZ
Nickel	0.014 ND	0.0040	mg/L	12/18/06	BSZ
Selenium	ND ND	0.010	mg/L	12/18/06	BSZ
Silver	ND ND	0.0020	mg/L	12/18/06	BSZ
			mg/L		
Thallium	ND	0.0050	mg/L	12/18/06	BSZ
Zinc	0.023	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW	ND	0.00020	mg/L	12/14/06	AM
(2) 841061208-28					
Date Collected: 12/8/2006 Matrix: Aqueous					
Trace Metals by 6010B					
Antimony	ND	0.010	mg/L	12/18/06	BSZ
Arsenic	ND	0.010	mg/L	12/18/06	BSZ
Barium	0.038	0.010	mg/L	12/18/06	BSZ
Beryllium	ND	0.0010	mg/L	12/18/06	BSZ
Cadmium	ND	0.0020	mg/L	12/18/06	BSZ
Chromium	ND	0.010	mg/L	12/18/06	BSZ
Copper	ND	0.010	mg/L	12/18/06	BSZ
Lead	0.012	0.0040	mg/L	12/18/06	BSZ
Nickel	ND	0.010	mg/L	12/18/06	BSZ
Selenium	ND	0.010	mg/L	12/18/06	BSZ
Silver	ND	0.0020	mg/L	12/18/06	BSZ
Thallium	ND	0.0050	mg/L	12/18/06	BSZ
Zinc	0.015	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW	ND	0.00020	mg/L	12/14/06	AM

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INORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, LLC
PL Report No: E612529
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/11/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(3) 841061208-29					
Date Collected: 12/8/2006 Matrix: Aqueous					
Trace Metals by 6010B	-				
Antimony	ND	0.010	mg/L	12/18/06	BSZ
Arsenic	ND	0.010	mg/L	12/18/06	BSZ
Barium	0.14	0.010	mg/L	12/18/06	BSZ
Beryllium	ND	0.0010	mg/L	12/18/06	BSZ
Cadmium	ND	0.0020	mg/L	12/18/06	BSZ
Chromium	ND	0.010	mg/L	12/18/06	BSZ
Copper	ND	0.010	mg/L	12/18/06	BSZ
Lead	ND	0.0040	mg/L	12/18/06	BSZ
Nickel	0.017	0.010	mg/L	12/18/06	BSZ
Selenium	ND	0.010	mg/L	12/18/06	BSZ
Silver	ND	0.0020	mg/L	12/18/06	BSZ
Thallium	ND	0.0050	mg/L	12/18/06	BSZ
Zinc	0.028	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW	ND	0.00020	mg/L	12/14/06	AM
(4) 841061208-30					
Date Collected: 12/8/2006 Matrix: Aqueous	<u> </u>				
Trace Metals by 6010B					
Antimony	ND	0.010	mg/L	12/18/06	BSZ
Arsenic	ND	0.010	mg/L	12/18/06	BSZ
Barium	0.15	0.010	mg/L	12/18/06	BSZ
Beryllium	ND	0.0010	mg/L	12/18/06	BSZ
Cadmium	ND	0.0020	mg/L	12/18/06	BSZ
Chromium	ND	0.010	mg/L	12/18/06	BSZ
Copper	0.015	0.010	mg/L	12/18/06	BSZ
Lead	ND	0.0040	mg/L	12/18/06	BSZ
Nickel	0.15	0.010	mg/L	12/18/06	BSZ
Selenium	ND	0.010	mg/L	12/18/06	BSZ
Silver	ND	0.0020	mg/L	12/18/06	BSZ
Thallium	ND	0.0050	mg/L	12/18/06	BSZ
Zinc	0.057	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW	ND	0.00020	mg/L	12/14/06	AM

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INORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, LLC
PL Report No: E612529
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/11/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter		Result	DL	Units	Completed	By Dilution
(5) 841061208-31						
Date Collected: 12/8/2006	Matrix: Aqueous					
Trace Metals by 6010B	matrix: riqueous					
Antimony		ND	0.010	mg/L	12/18/06	BSZ
Arsenic		ND	0.010	mg/L	12/18/06	BSZ
Barium		0.83	0.010	mg/L	12/18/06	BSZ
Beryllium		0.0018	0.0010	mg/L	12/18/06	BSZ
Cadmium		0.0034	0.0020	mg/L	12/18/06	BSZ
Chromium		0.092	0.010	mg/L	12/18/06	BSZ
Copper		0.073	0.010	mg/L	12/18/06	BSZ
Lead		1.9	0.0040	mg/L	12/18/06	BSZ
Nickel		0.12	0.010	mg/L	12/18/06	BSZ
Selenium		ND	0.010	mg/L	12/18/06	BSZ
Silver		ND	0.0020	mg/L	12/18/06	BSZ
Thallium		ND	0.0050	mg/L	12/18/06	BSZ
Zinc		0.73	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW		ND	0.00020	mg/L	12/14/06	AM
(6) 841061208-32						
Date Collected: 12/8/2006	Matrix: Aqueous					
Trace Metals by 6010B						
Antimony		ND	0.010	mg/L	12/18/06	BSZ
Arsenic		ND	0.010	mg/L	12/18/06	BSZ
Barium		0.21	0.010	mg/L	12/18/06	BSZ
Beryllium		0.0087	0.0010	mg/L	12/18/06	BSZ
Cadmium		ND	0.0020	mg/L	12/18/06	BSZ
Chromium		0.036	0.010	mg/L	12/18/06	BSZ
Copper		0.018	0.010	mg/L	12/18/06	BSZ
Lead		0.098	0.0040	mg/L	12/18/06	BSZ
Nickel		0.054	0.010	mg/L	12/18/06	BSZ
Selenium		ND	0.010	mg/L	12/18/06	BSZ
Silver		ND	0.0020	mg/L	12/18/06	BSZ
Thallium		ND	0.0050	mg/L	12/18/06	BSZ
Zinc		0.17	0.010	mg/L	12/18/06	BSZ
Mercury by SW-846 7470A in GW		ND	0.00020	mg/L	12/14/06	AM

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061208-27
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/12/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	72	70%-130%	
2,5-dibromotoluene #2	72	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 1

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-27

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32643.D

Forestand Acetone	CAS No.	Parameter	Result	DL
108.86-1 Bromobenzene	67-64-1	Acetone	ND	10
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 1.0 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 108-90-7 Chlorocthane ND 1.0 108-90-7 Chloroform ND 1.0 74-87-3 Chloroform ND 1.0 14-87-3 Chloroformethane ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-93-4<	71-43-2	Benzene	ND	1.0
75-27-4 Bromoform ND 1.0 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorobenzene ND 1.0 76-66-3 Chloroform ND 1.0 74-87-3 Chlorobenzene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 19-2-8	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-34 1.2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromomethane ND 0.50	74-97-5	Bromochloromethane	ND	1.0
74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorotelhane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotelhane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-joromo-chloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50	75-27-4	Bromodichloromethane	ND	1.0
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Burylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorobenzene ND 1.0 67-66-3 Chloroform ND 1.0 67-48-8 2-Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 104-48-1 Dibromochlane (EDB) ND 0.50 14-95-3 Dibromochlane (EDB) ND 0.50 106-93-4 1,2-Dichlorobenzene ND 1.0	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorochtane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 104-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 105-93-4 1,2-Dibromo-shane ND 0.50 40-93-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND	104-51-8	· · · · · · · · · · · · · · · · · · ·	ND	1.0
98-06-6 tert-Buylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloromethane ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.5 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 14-95-3 Dibromochloromethane ND 0.50 106-93-4 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND<		· · · · · · · · · · · · · · · · · · ·		
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 104-48-1 Dibromochloromethane ND 0.50 104-48-1 Dibromochloromethane (EDB) ND 0.50 104-48-1 Dibromomethane ND 0.50 105-50-3 1,2-Dichlorobenzene ND 1.0 105-50-3 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 105-46-7 1,4-Dichlorobenzene ND				
56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorotethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochlane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 10-6-20 1,2-Dichloroethane ND </td <td></td> <td>· · · · · · · · · · · · · · · · · · ·</td> <td>ND</td> <td></td>		· · · · · · · · · · · · · · · · · · ·	ND	
108-90-7 Chloroebnezene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 14-95-3 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorodifluoromethane				
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,4-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 175-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane <td< td=""><td></td><td></td><td></td><td></td></td<>				
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromoethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 55-71-8 Dichlorodifluoromethane ND 1.0 75-71-8 Dichlorodhene ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-69-2 cis-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND		Chloroethane		
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isporpyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 57-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorodethane ND 1.0 107-06-2 1,2-Dichlorotethane ND 1.0 156-59-2 cis-1,2-Dichlorotethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichlorop		Chloroform		
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60-29-7 Diethyl ether ND 1.0				
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			ND	

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612529 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 (continued) Sample Description: 841061208-27

Date Collected: 12/8/2006 Matrix: Aqueous

Date Received: 12/11/2006 Percent Moisture: N/A

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/14/06 By: GP Dilution Factor: 1

Method: 8260B

Soil Extract Volume:

Method: 8260B Soil Extract Volume:
QC Batch#: 51242 Lab Data File: M32643.D

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lir	nits	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	107%	91%-109%
Bromofluorobenzene	90%	87%-105%
Toluene-d8	104%	92%-105%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061208-28
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/13/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	78	70%-130%	
2,5-dibromotoluene #2	85	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 2

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-28

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32644.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	1.0
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
108-20-3	Di-isopropyl ether (DIPE)	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612529 PL Sample No: 2 (continued) Sample Description: 841061208-28

Date Collected: 12/8/2006 Date Received: 12/11/2006 Date Extracted: By: Date Analyzed: 12/14/06 By: GP

Dilution Factor: 1 Method: 8260B Soil Extract Volume: Lab Data File: M32644.D QC Batch#: 51242

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	108%	91%-109%	
Bromofluorobenzene	88%	87%-105%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	108%	91%-109%
Bromofluorobenzene	88%	87%-105%
Toluene-d8	103%	92%-105%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061208-29
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/13/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	78	70%-130%
2,5-dibromotoluene #2	85	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	2.1	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Custom

PL Report No: E612529 PL Sample No: 3

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/13/06 By: GP

Method: 8260B QC Batch#: 51243 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-29

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume:

Lab Data File: J28736.D;M32648.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	1.0
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
108-20-3	Di-isopropyl ether (DIPE)	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	8.0	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612529 PL Sample No: 3 (continued) Sample Description: 841061208-29

Date Collected: 12/8/2006 Matrix: Aqueous Date Received: 12/11/2006 Percent Moisture: N/A Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/13/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume:

QC Batch#: 51243 Lab Data File: J28736.D;M32648.D

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	1.8	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	240	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	1.8	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	150	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	100%	91%-109%	
Bromofluorobenzene	89%	87%-105%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	100%	91%-109%
Bromofluorobenzene	89%	87%-105%
Toluene-d8	103%	92%-105%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061208-30
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/13/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	75	70%-130%
2,5-dibromotoluene #2	81	70%-130%

TARGETED VPH ANALYTES

Analyte	Results QL		Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 4

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-30

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32645.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	1.0
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
108-20-3	Di-isopropyl ether (DIPE)	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612529 PL Sample No: 4 (continued) Sample Description: 841061208-30

Date Collected: 12/8/2006 Date Received: 12/11/2006 Date Extracted: By: Date Analyzed: 12/14/06 By: GP

Sample Weight/Volume: Dilution Factor: 1 Method: 8260B Soil Extract Volume: QC Batch#: 51242 Lab Data File: M32645.D

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	6.6	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	6.6	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	108%	91%-109%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	108%	91%-109%
Bromofluorobenzene	89%	87%-105%
Toluene-d8	101%	92%-105%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061208-31
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/13/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	79	70%-130%
2,5-dibromotoluene #2	84	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 5

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-31

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32646.D

Forestand Acetone	CAS No.	Parameter	Result	DL
108.86-1 Bromobenzene	67-64-1	Acetone	ND	10
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 1.0 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 108-90-7 Chlorocthane ND 1.0 108-90-7 Chloroform ND 1.0 74-87-3 Chloroform ND 1.0 14-87-3 Chloroformethane ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-93-4<	71-43-2	Benzene	ND	1.0
75-27-4 Bromoform ND 1.0 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorobenzene ND 1.0 76-66-3 Chloroform ND 1.0 74-87-3 Chlorobenzene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 19-2-8	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-34 1.2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromomethane ND 0.50	74-97-5	Bromochloromethane	ND	1.0
74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorotelhane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotelhane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 108-20-3 Di-joromo-chloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50	75-27-4	Bromodichloromethane	ND	1.0
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Burylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorobenzene ND 1.0 67-66-3 Chloroform ND 1.0 67-48-8 2-Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 104-48-1 Dibromochlane (EDB) ND 0.50 14-95-3 Dibromochlane (EDB) ND 0.50 106-93-4 1,2-Dichlorobenzene ND 1.0	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorochtane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 108-20-3 Di-isopropyl ether (DIPE) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 104-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 105-93-4 1,2-Dibromo-shane ND 0.50 40-93-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND	104-51-8	· · · · · · · · · · · · · · · · · · ·	ND	1.0
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60-29-7 Diethyl ether ND 1.0				
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			ND	

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612529 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 (continued) Sample Description: 841061208-31

Date Collected:12/8/2006Matrix: AqueousDate Received:12/11/2006Percent Moisture: N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:12/14/06By: GPDilution Factor: 1Method:8260BSoil Extract Volume:

Method: 8260B Soil Extract Volume:
QC Batch#: 51242 Lab Data File: M32646.D

Units: ug/L

CAS No.	o. Parameter			DL
	Ethyl tertiary-butyl ether (EtBE)		ND	1.0
100-41-4	Ethylbenzene		ND	1.0
87-68-3	Hexachlorobutadiene		ND	0.50
591-78-6	2-Hexanone		ND	5.0
98-82-8	Isopropylbenzene		ND	1.0
99-87-6	4-Isopropyltoluene		ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	5.0
75-09-2	Methylene chloride		ND	5.0
91-20-3	Naphthalene		ND	1.0
103-65-1	n-Propylbenzene		ND	1.0
100-42-5	Styrene		ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)		ND	1.0
109-99-9	Tetrahydrofuran		ND	1.0
96-18-4	1,2,3-Trichloropropane		ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane		ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane		ND	0.50
127-18-4	Tetrachloroethene (PCE)	Tetrachloroethene (PCE)		1.0
108-88-3	Toluene	Toluene		1.0
87-61-6	1,2,3-Trichlorobenzene	1,2,3-Trichlorobenzene		1.0
120-82-1	1,2,4-Trichlorobenzene	1,2,4-Trichlorobenzene		1.0
71-55-6	1,1,1-Trichloroethane	1,1,1-Trichloroethane		1.0
79-00-5	1,1,2-Trichloroethane		ND	1.0
79-01-6	Trichloroethene (TCE)		ND	1.0
75-69-4	Trichlorofluoromethane		ND	1.0
95-63-6	1,2,4-Trimethylbenzene		ND	1.0
108-67-8	1,3,5-Trimethylbenzene		ND	1.0
75-01-4	Vinyl chloride	Vinyl chloride		1.0
95-47-6	o-Xylene	o-Xylene		1.0
	m,p-Xylenes		ND	1.0
Surrogate	Recovery	Limits		
1,2-Dichloroethane-d4	108%	91%-109%		
Bromofluorobenzene	89%	87%-105%		
Toluene-d8	103%	92%-105%		

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061208-32
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Analyzed:	12/13/06	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	75	70%-130%
2,5-dibromotoluene #2	79	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 6

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-32

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32647.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	1.0
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
108-20-3	Di-isopropyl ether (DIPE)	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Customer: Fuss & O'Neill Laboratory: Premier Laboratory, LLC Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612529 PL Sample No: 6 (continued) Sample Description: 841061208-32

Date Collected: 12/8/2006 Date Received: 12/11/2006 Date Extracted: Date Analyzed: 12/14/06 By: GP

Sample Weight/Volume: Dilution Factor: 1 Method: 8260B Soil Extract Volume: QC Batch#: 51242 Lab Data File: M32647.D

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	43	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	40	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	108%	91%-109%	
Bromofluorobenzene	89%	87%-105%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	108%	91%-109%
Bromofluorobenzene	89%	87%-105%
Toluene-d8	101%	92%-105%

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Laboratory: Premier Laboratory, LLC

PL Report No: E612529 PL Sample No: 7

Date Collected: 12/8/2006

Date Received: 12/11/2006

Date Extracted: By:

Date Analyzed: 12/14/06 By: GP

Method: 8260B QC Batch#: 51242 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841061208-33

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: M32639.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	1.0
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
108-20-3	Di-isopropyl ether (DIPE)	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612529 PL Sample No: 7 (continued) Sample Description: 841061208-33

Date Collected: 12/8/2006 Date Received: 12/11/2006 Date Extracted: By: Date Analyzed: 12/14/06 By: GP

Sample Weight/Volume: Dilution Factor: 1 Method: 8260B Soil Extract Volume: QC Batch#: 51242 Lab Data File: M32639.D

Units: ug/L

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1.0
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
1748-03-8	Tertiary-amyl methyl ether (TAME)	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
	m,p-Xylenes	ND	1.0
Surrogate	Recovery I	Limits	
1,2-Dichloroethane-d4	107%	91%-109%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	107%	91%-109%
Bromofluorobenzene	91%	87%-105%
Toluene-d8	104%	92%-105%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061208-27
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/13/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	76	40%-140%
2-Bromonaphthalene	101	40%-140%
2-Fluorobiphenyl	95	40%-140%
o-Terphenyl	72	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061208-28
Preservative	HCL	•	
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/13/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	73	40%-140%
2-Bromonaphthalene	100	40%-140%
2-Fluorobiphenyl	94	40%-140%
o-Terphenyl	72	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061208-29
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/12/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	70	40%-140%
2-Bromonaphthalene	96	40%-140%
2-Fluorobiphenyl	91	40%-140%
o-Terphenyl	69	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061208-30
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/13/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	51	40%-140%
2-Bromonaphthalene	68	40%-140%
2-Fluorobiphenyl	70	40%-140%
o-Terphenyl	45	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061208-31
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/13/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	50	40%-140%
2-Bromonaphthalene	72	40%-140%
2-Fluorobiphenyl	77	40%-140%
o-Terphenyl	51	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612529	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061208-32
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	12/8/2006		
Date Received:	12/11/2006	Matrix:	Aqueous
Date Extracted:	12/12/06	Percent Moisture:	N/A
Date Analyzed:	12/13/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	55	40%-140%
2-Bromonaphthalene	100	40%-140%
2-Fluorobiphenyl	95	40%-140%
o-Terphenyl	72	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

Page 32 of 32



□ 146 Hartford Road, Manchester, CT 06040

56 Quarry Road, Trumbull, CT 06611

□ 1419 Richland Street, Columbia, SC 29201

□ 78 Interstate Drive, West Springfield, MA 01089

□ 610 Lynndale Court, Suite E, Greenville, NC 27858

□ 24 Madison Avenue Extension, Albany, NY 12203

10636

275 Promenade Street, Suite 350, Providence, RI 0290

□ 80 Washington Street, Suite 301, Poughkeepsie, NY 12601

Turnaround

□ Other_

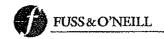
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PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

		YES NO COMMENTS
1. 2. 3.	SDG Project Narratives Traffic Report Volatiles Data	
	a. Sample Data Target Compound List (TCL) Results Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library matches Percent solids calculations	
	b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary	
	c. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Duplicate Data	X = For 8260 X = analysis only
4.	Semivolatiles Data a. QC Summary Surrogate Percent Recovery Summary MS/MSD Summary	X
	Method Blank Summary Tuning and Mass Calibration	



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

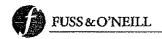
		YES	NO.	COMMENT
b.	Sample Data TCL Results	Ж		
	Tentatively Identified Compounds		` \(\overline{\sqrt{2}}\)	
	Reconstructed total ion chromatograms (RIC) for each			
	Sample	攵		
	For each sample:		`	
	<u> </u>			
			×.	
			X	
	GPC chromatograms (if GPC performed)		Ο.	<i>N</i> /4
c.	Standards Data (all instruments)			•
	Initial Calibration Data	X		
	RICs and Quan Reports for all Standards	X		
	Continuing Calibration)B,		
	RICs and Quan Reports for all Standards	X		
	Internal Standard Areas Summary			U/A
	Internal Standard Areas Summary			
đ.	Raw QC Data			A 5
				NA
	Blank Data	X		
	Matrix Spike Data	(D))[2]	
	I'CL Results Tentatively Identified Compounds Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass pectra of TCL compounds Mass spectra of TICs with 3 best library matches GPC chromatograms (if GPC performed) Standards Data (all instruments) Initial Calibration Data LICs and Quan Reports for all Standards Continuing Calibration LICs and Quan Reports for all Standards Internal Standard Areas Summary Internal Standards Internal		[X] _	······································
Mis	cellaneous Data			
	Original preparation and analysis forms or copies of preparation	,		
	and analysis log book pages	X		
	Internal sample & sample extract transfer chain-of custody	, .		
3	records		× _	····
	Screening Records)	
4	All instrument output, including strip charts from screening		X _	·
	activities (describe or list)			



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

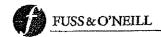
				YES NO	<u>COMMENTS</u>	
6.		ustody Records Sample Log-in Sheet (Lab & DC1) Miscellaneous Shipping/Receiving Rec	ords (describe or list)	X 0 - X 0	NA	
7.	Internal Lal	Sample Transfer Records and Track Sheets (describe or list) Somyle translo	9	X ==		
8.	Other Reco	rds (describe or list) INSTRUMENT REN	logBilot	X		
9.	Comments:	* See laboratory Quality Assurance	Plan for limits.			
	Complete (Lab)	4.J. AL (1)	Michae / I (Printed Name/Title)	Y lall	um 1/4/02 Date	
		nat the above information is true and according analyses will be made available for review				ì
	Certified l (Lab)		Mober Stevens (Printed Name/Title)	as Lab Dicc	dor 1-15-07 Date	



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

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SDG Project Narratives	Ø]		
Inorganic Analysis Data Sheet	₫]		
Initial and Continuing Calibration Verification	12]		
CRDL Standard for AA and ICP	127]		
Blanks	团	Е]		
ICP Interference Check Sample	12]		
Spike Sample Recovery	M]		
Post Digest Spike Sample Recovery		V		NA	
Duplicates	5 Z1			·	
Laboratory Control Sample	12				
Standard Addition Results		Z	į		
ICP Serial Dilutions	瓦				
Instrument Detection Limits, Quarterly	囟				
ICP Interelement Correction Factors, Annually	12]				
ICP Linear Ranges Quarterly	IZ)				
Preparation Log	[☑]				
Analysis Run Log	M				
ICP Raw Data					
Furnace AA Raw Data		Ø		N ↑	
· Mercury Raw Data	◩				
Percent Solids Calculations			•		
Digestion Logs	Ì ∑Ì				
EPA Shipping/Receiving Records			•		
(List all individual records)			_		
Chain-of Custody Records	121				
Sample Log-In sheet					
Miscellaneous Shipping/Receiving Records			_		
(List all individual records)			-		•



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS . (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

YES NO COMMENTS

25.	Internal Lab Sample Transfer Records a (Describe or List)	nd Tracking Sheets	_	
26.	Internal Original Sample Preparation and (Describe or List Preparation Records Analysis Records Description	d analysis Records		[]
27.	Other Records (Describe or List)		,	
28.	Comments:			
* *	See laboratory Quality Assurance Plan fo	or limits.		
Comple Lab)	sted by: B. Suppenki (Signature)	Barbaw-Słupryczyni (Printed Name/Title)	ti / Pro	duction Hamag. 1/8/07 Date
ssociat	that the above information is true and aced with the above analyses will be made a tion of this document.			
Certified Lab)	1 by: B. Auph Cyntki B. (Signature)	avbaw Suprycynski Printed Name/Title)	Produc	tion Hang. 1807 Date

Spike Recovery and RPD Summary Report - WATER

Method : C:\HPCHEM\1\METHODS\8260BDIR.M

Title : EPA METHOD 8260B

Last Update : Wed Jan 03 12:23:29 2007

Response via : Initial Calibration

Non-Spiked Sample: J28732.D

Spike Spike

Sample Duplicate Sample

File ID : J28729.D J28730.D

Sample : VLCS1213 DUP

Acq Time: 13 Dec 06 12:23 pm 13 Dec 06 12:50 pm

Compound	Sample Conc	Spike Added	Spike Res	Dup Res	Spike %Rec	Dup %Rec	RPD	QC RPD	Limits % Rec
Dichlorodifluorometh	0.0	50	58	54	116	107	8	25	70-130
Chloromethane	0.0	50	42	39	84	78	7	25	70-130
Vinyl chloride	0.0	50	47	43	94	87	8	25	70-130
Bromomethane	0.0	50	47	44	100	93	7	25	70-130
Chloroethane	0.0	50	50	48	113	108	5	25	70-130
Trichlorofluorometha	0.0	50	41	37	81	74	9	25	70-130
Acetone	0.0	50	46	47	92	93	1	25	70-125
1,1-Dichloroethene	0.0	50	57	53	114	106	7	25	74-130
Carbon disulfide	0.0	50	42	39	84	78	7	25	70-130
Methylene chloride	0.0	50	55	53	116	112	4	25	70-127
Methyl-tertbutyl eth	0.0	50	50	49	100	98	2	25	80-117
trans-1,2-Dichloroet	0.0	50	54	51	108	103	5	25	73-123
Tertiary-butyl alcoh	0.0	500	520	514	104	103	1	25	70-130
Di-isopropyl ether (0.0	50	50	49	100	98	2	25	70-130
1,1-Dichloroethane	0.0	50	48	46	96	93	4	25	75-121
Ethyl tertiary-butyl	0.0	50	50	49	100	99	1	25	70-130
2,2-Dichloropropane	0.0	50	64	64	128	128	1	25	70-130
cis-1,2-Dichloroethe	0.0	50	52	50	104	100	4	25	78-117
Chloroform	0.0	50	48	46	96	92	4	25	72-118
Bromochloromethane	0.0	50	50	49	100	98	2	25	70-130
tetrahydrofuran	0.0	50	43	43	87	8.7	0	25	70-130
1,1,1-Trichloroethan	0:0	50	52	49.	103	98	6	25	75-124
1,1-Dichloropropene	0.0	50	53	50	105	101	4	25	81-118
Carbon tetrachloride	0.0	50	52	49	104	99	5	25	74-127
Benzene	0.0	50	49	46	98	93	5	25	76-118
Tertiary-amyl methyl	0.0	50	51	50	101	100	1	25	70-130
1,2-Dichloroethane	0.0	50	46	45	91	90	1	25	70-114
Trichloroethene	0.0	50	48	46	97	92	5	25	74-117
1,2-Dichloropropane	0.0	50	47	45	94	90	4	25	73-118
Bromodichloromethane	0.0	50	54	52	109	104	4	25	73-122
Dibromomethane	0.0	50	49	48	98	96	1	25	70-130
1,4-Dioxane	0.0	50	56	53	113	106	6	25	70-130
cis-1,3-Dichloroprop	0.0	50	49	49	98	99	1	25	70-120
Toluene	0.0	50	49	47	98	95	4	25	76-117
trans-1,3-Dichloropr	0.0	50	46	46	93	92	1	25	70-119
1,1,2-Trichloroethan	0.0	50	47	47	94	95	1	25	71-115
1,2-Dibromoethane	0.0	50	48	49	97	98	1	25	86-107
1,3-Dichloropropane	0.0	50	49	49	98	98	0	25	79-114
Tetrachloroethene	0.0	50	47	46	94	91	3	25	74-120
Chlorobenzene	0.0	50	48	47	97	94	3	25	76-119
1,1,1,2-Tetrachloroe	0.0	50	48	46	96	92	4	25	75-116
Ethylbenzene	0.0	50	52	49	103	99	5	25	78-122
o-Xylene	0.0	50	49	47	97	95	3	25	81-116
4	• •		ļ	'		. !		,	1

Styrene	0.0	50	53	52	105	103	2	25	81-120
Bromoform	0.0	50	52	51	104	103	1	25	70-122
Isopropylbenzene	0.0	50	52	50	113	108	4	25	82-119
1,1,2,2-Tetrachloroe	0.0	50	54	54	119	119	1	25	70-123
n-Propylbenzene	0.0	50	52	50	103	100	3	25	83-126
Bromobenzene	0.0	50	52	50	112	109	3	25	70-130
2-Chlorotoluene	0.0	50	52	50	111	108	3	25	79-121
4-Chlorotoluene	0.0	50	45	44	89	88	1	25	78-119
1,3,5-Trimethylbenze	0.0	50	4.8	47	95	94	1	25	83-123
tert-Butylbenzene	0.0	50	41	41	83	83	0	25	81-119
1,2,4-Trimethylbenze	0.0	50	51	49	102	99	3	25	79-121
sec-Butylbenzene	0.0	50	48	46	96	91	5	25	70-130
1,3-Dichlorobenzene	0.0	50	48	47	96	94	1	25	73-120
1,4-Dichlorobenzene	0.0	50	47	46	94	92	3	25	74-121
n-Butylbenzene	0.0	50	42	40	84	80	6	25	70-129
1,2-Dichlorobenzene	0.0	50	48	48	97	96	1	25	74-121
1,2,4-Trichlorobenze	0.0	5.0	41	41	82	82	2	25	70-124
Hexachlorobutadiene	0.0	50	41	41	82	81	1	25	70-122
Naphthalene	0.0	50	42	42	84	84	2	25	70-128
1,2,3-Trichlorobenze	0.0	50	41	40	82	80	1	25	70-130

8260BDIR.M

Wed Jan 03 13:43:00 2007

MS8

Spike Recovery and RPD Summary Report - WATER

Method : C:\HPCHEM\1\METHODS\8260BMCP.M (RTE Integrator)

Title

Last Update : Wed Jan 03 13:38:48 2007

Response via: Initial Calibration

Non-Spiked Sample: M32638.D

Spike Spike

Sample Duplicate Sample

File ID: M32634.D | M32635.D

Sample: VLCS1214 | VLCS1214 DUP
Acq Time: 14 Dec 2006 9:22 am | 14 Dec 2006 9:53 am

Compound	Sample Conc	Spike Added	Spike Res	Dup Res	Spike %Rec	Dup %Rec	RPD	QC RPD	Limits % Rec
Dichlorodifluorometh	1 0.0	50	62	64	124	128	3	1 25	70-130
Chloromethane	0.0	50	42	42	83	85	2	25	70-130
Vinyl chloride	0.0	50	49	50	98	100	1 2	25	70-130
Bromomethane	0.1	50	53	54	1 107	108	1 1	25	70-130
Chloroethane	0.0	50	62	62	124	124	i	25	70-130
Trichlorofluorometha		50	62	62	123	125	2	25	70-130
	0.0	50	42	43	85	86	2	25	83-124
Diethyl ether	1 0.0	i 50	42	48	88	1 96	9	25	70-125
Acetone	0.0	50	59	59	1 117	1118	1 1	25	74-130
1,1-Dichloroethene		50	45	44	89	88	1	25	70-130
Carbon disulfide	0.0	50	55.	58	1 111	116	5	25	70-127
Methylene chloride	1 0.0	50	50	53	1 100	1 106	1 6	25	80-117
Methyl-tertbutyl eth		50	1 53 I	54	1 106	100	3	1 25	73-123
trans-1,2-Dichloroet		•	33	502	89	100	1 12	25	70-130
Tertiary-butyl alcoh		500	51	52	102		3	25	70-1301
Di-isopropyl ether (50	51	52	102	103	2	25	75-121
1,1-Dichloroethane	0.0	50 50	55	56	102	112	2	25	70-130
Ethyl tertiary-butyl		•			1 109	1113	3	25	70-130
2,2-Dichloropropane	1 0.0	50 50	55 54		1 107	113	3	25	78-117
cis-1,2-Dichloroethe		•	51	52	107	105	1 2	25	72-118
Chloroform	0.0	50	31	49	95	99	1 4	25	70-130
Bromochloromethane	0.0	50		40	78	81	3	1 25	70-130
Tetrahydrofuran	0.0	50	39 56	58	1112	1115	1 3	1 25	75-124
1,1,1-Trichloroethan		50	30	50	98	100	2	25	81-118
1,1-Dichloropropene	0.0	50 50	54	54	1 108	107	1 1	25	74-127
Carbon tetrachloride		50	51	51	100	103	1 1	25	76-118
Benzene	0.0	, 50 I 50	53	56	1 107	1112	5	25	70-130
Tertiary-amyl methyl	1 0.2	50	54	55	108	110	1	25	70-114
1,2-Dichloroethane	1 0.0	50 50	49	49	1 98	99	1	25	74-117
Trichloroethene	0.0	50 50	52	55	104	110	5	25	73-118
1,2-Dichloropropane	0.0	50	52	54	1 104	108	4	25	70-130
Dibromomethane	1	50	57	60	1114	1119	, <u>,</u> ,	25	73-122
Bromodichloromethane	0.0	50	39	43	79	86	i 8	25	70-130
1,4-Dioxane	:	50	53	51	106	102	4	25	70-120
cis-1,3-Dichloroprop Toluene	0.0	50	55	53	109	107	2	25	76-117
trans-1,3-Dichloropr	•	50	51	49	102	97	5	25	70-119
1,1,2-Trichloroethan		50	54	52	107	105	2	25	71-115
1,2-Dibromoethane	0.0	50	50	50	100	101	1	25	86-107
1,3-Dichloropropane	0.0	50	55	56	109		3	25	79-114
Tetrachloroethene	0.0		53	51	107		4	25	74-120
Chlorobenzene	0.0	50	53	53	106	105	1	25	76-119
1,1,1,2-Tetrachloroe	•	50	54	54	108	109	j 1	25	75-116
Ethylbenzene	0.0	50	59	58	117	116	1	25	78-122
m+p-Xylenes	0.0	100	118	114	118	114	4	25	80-124
	0.0	50	55	53	110	107	-	25	81-116
o-Xylene	0.0		56	54	113	109	4	25	81-120
Styrene		50	51	51	103	103	0	25	70-1221
Bromoform	0.0	50	54	52	108	105	j 3	25	82-119
Isopropylbenzene 1,1,2,2-Tetrachloroe		50	49	51	99	102	j 3	25	70-123
n-Propylbenzene	0.0	50	58	57	116	114	j 2	25	83-126
n-propyrdenzene Bromobenzene	0.0	50	53	54	107	109	j 2	25	70-130
2-Chlorotoluene	0.1			50	103	100	j 3	25	79-121
Z CHIOLOCOLUCIO	, ,,,,			-					

4-Chlorotoluene 1,3,5-Trimethylbenze tert-Butylbenzene 1,2,4-Trimethylbenze sec-Butylbenzene 4-Isopropyltoluene 1,3-Dichlorobenzene 1,4-Dichlorobenzene n-Butylbenzene 1,2-Dichlorobenzene	0.0	50 50 50 50 50 50 50 50	56 55 52 57 52 53 55 50 48 53	54 112 53 111 52 104 55 113 49 104 51 106 55 110 48 99 47 97 52 105	108 107 104 111 99 101 109 95 93	4 25 4 25 0 25 2 25 5 25 5 25 0 25 4 25 4 25 1 25	78-119 83-123 81-119 79-121 70-130 70-130 73-120 74-121 70-129 74-121
1,4-Dichlorobenzene n-Butylbenzene	0.1 0.0 0.0	50 50	50 48	48 99 47 97	95	4 25	74-121 70-129 74-121 70-124 70-122 70-128

- Fails Limit Check

8260BMCP.M Wed Jan 03 13:39:19 2007

FORM 3 WATER 8260B MATRIX SPIKE/MATRIX SPIKE DUPLICATE RECOVERY

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/14/06

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Sample No.: E612529-1C

Location: Franklin, MA

	COMPOUND	SPIKE ADDED (ug/L)	SAMPLE CONCENTRATION (ug/L)	MS CONCENTRATION (ug/L)	MS % REC #	QC LIMITS REC
1:		=======	_=========	=======================================	======	=====
	1,1-Dichloroethene	50.00	0	61.42	123	34-176
	Benzene	50.00	0	52.21	104	69-115
1	Chlorobenzene	50.00	0	51.40	103	78-108
	Toluene	50.00	0	53.43	107	75-109
	Trichloroethene (TCE)	50.00	0	50.13	100	70-116
1_				· 		ii

	SPIKE ADDED	MSD CONCENTRATION	MSD	 %	OC LIMITS
COMPOUND	(ug/L)	(ug/L)	REC #	,	. ~
			=====	=====	=======================================
1,1-Dichloroethene	50.00	67.08	134	8.56	71 34-176
Benzene	50.00	53.24	106	1.90	23 69-115
Chlorobenzene	50.00	50.78	102	0.98	15 78-108
Toluene	50.00	54.17	108	0.93	17 75-109
Trichloroethene (TCE)	50.00	50.54	101	1.0	23 70-116
				·	

Column to be used to flag recovery and RPD values with an asterisk

RPD: 0 out of 5 outside limits

Spike Recovery: 0 out of 10 outside limits

COMMENTS:

FILE: M32643.D

^{*} Values outside of QC limits

FORM 4 8260B METHOD BLANK SUMMARY

 VBLK1214
VDDRIZZII

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Lab File ID: M32638.D

Lab Sample ID: VBLK1214

Matrix: (soil/water) Water Date Analyzed: 12/14/06

Instrument ID: MS11

Date Extracted:

Time Analyzed: 1126

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	LAB	CLIENT	LAB	DATE
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED
	_=========			========
		•	M32643.D	12/14/06
		•	M32644.D	12/14/06
			M32645.D	12/14/06
	•		M32646.D	12/14/06
		•	M32647.D	12/14/06
		•	M32639.D	12/14/06
		•	M32634.D	12/14/06
	E612529-1CMS		M32650.D	12/14/06
		E612529-1CMSD	M32651.D	12/14/06
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27 28]	
291				I
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30				

COMMENTS:	

FORM 4 8260B METHOD BLANK SUMMARY

VBLK1213	

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Lab File ID: J28732.D

Lab Sample ID: VBLK1213

Matrix: (soil/water) Water

Date Analyzed: 12/13/06

Instrument ID: MS8

Date Extracted:

Time Analyzed: 1346

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	LAB	CLIENT	LAB	DATE
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED
	1			
	-	•	J28736.D;M3264	
	•	VLCS1213	J28729.D	12/13/06
03	·			
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05	·			
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COMMENTS:	
	5

Data File : D:\HPCHEM\1\DATA\061214M.B\M32638.D

Vial: 6

Acq On : 14 Dec 2006 11:26 am : VBLK1214

Operator: GP

Sample Misc : 8260B()

: MS11 Inst Multiplr: 1.00

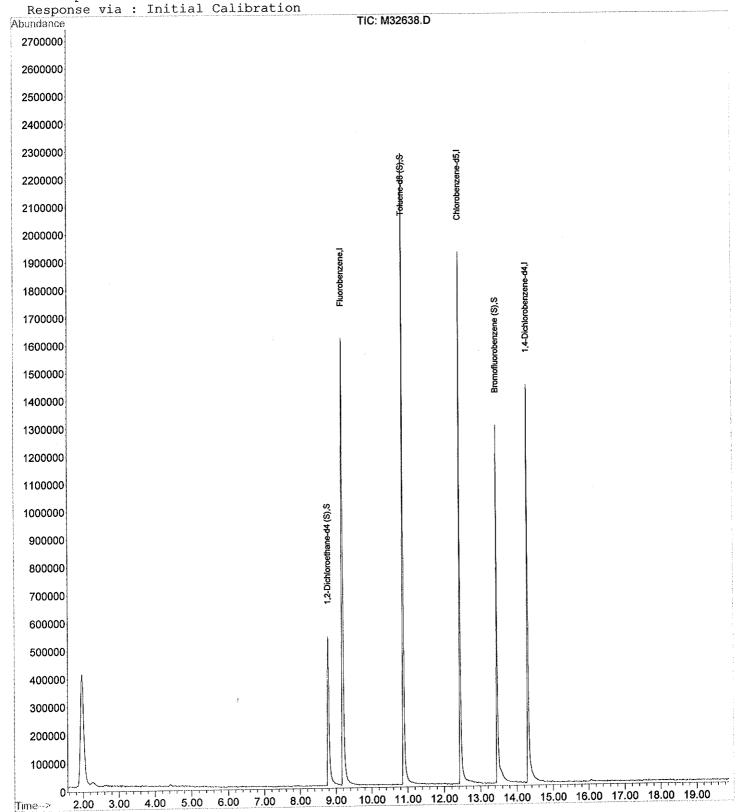
MS Integration Params: rteint.p

Quant Results File: 8260BMCP.RES Quant Time: Dec 14 11:49 2006

: C:\HPCHEM\1\METHODS\8260BMCP.M (RTE Integrator) Method

Title

Last Update : Wed Jan 03 12:15:52 2007



(QT Reviewed)

Data File : D:\HPCHEM\1\DATA\061214M.B\M32638.D

Quant Time: Dec 14 11:49 2006

Operator: GP Inst : MS11

Quant Results File: 8260BMCP.RES

Vial: 6

Acq On : 14 Dec 2006 11:26 am Sample : VBLK1214 Misc : 8260B() Multiplr: 1.00 MS Integration Params: rteint.p

Quant Method: C:\HPCHEM\1\METHODS\8260BMCP.M (RTE Integrator)

Title

Last Update : Thu Dec 14 10:34:26 2006 Response via : Initial Calibration

DataAcq Meth: 8260BMCP

Internal Standards	R.T.	QIon	Response	Conc Ur	nits Dev	(Min)
1) Fluorobenzene 43) Chlorobenzene-d5 63) 1,4-Dichlorobenzene-d4	9.19 12.43 14.31	96 117 152	1910585 995851 466920	50.00 50.00 50.00	ppb	0.00 0.00 0.00
System Monitoring Compounds 34) 1,2-Dichloroethane-d4 (S) Spiked Amount 50.000	8.78	65	558761m Recove	50.70	ppb 101.40%	0.00
46) Toluene-d8 (S)	10.87	98	1563709	51.49		0.00
Spiked Amount 50.000 52) Bromofluorobenzene (S) Spiked Amount 50.000	13.45	176	Recove 402131 Recove	45.88	102.98% ppb 91.76%	0.00

Target Compounds

Qvalue

Data File : C:\HPCHEM\1\DATA\061213J.B\J28732.D

Acq On : 13 Dec 06 1:46 pm

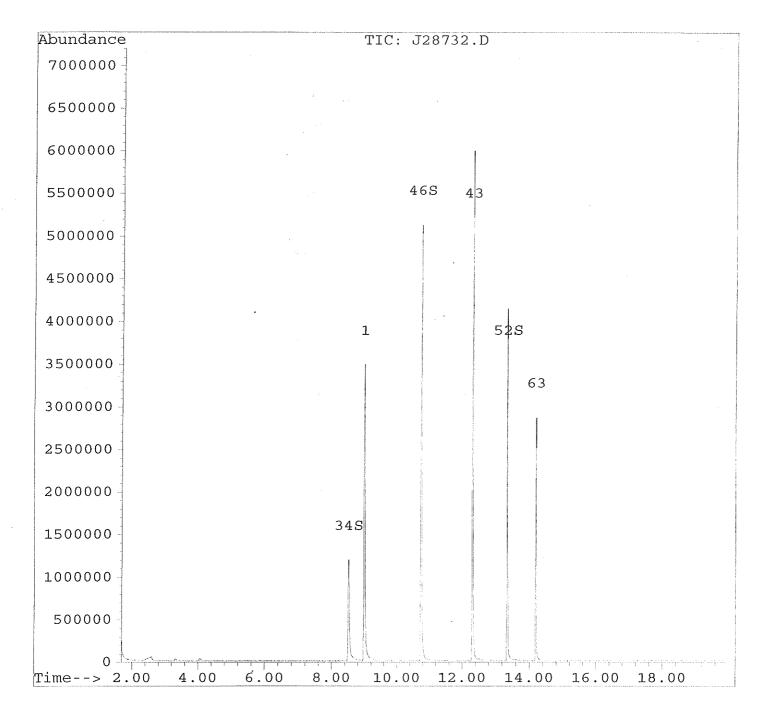
Sample : VBLK1213

Misc : 8260B() Quant Time: Jan 3 13:27 2007

Method : C:\HPCHEM\1\METHODS\8260BDIR.M

: EPA METHOD 8260B Title

Last Update : Wed Jan 03 12:23:29 2007 Response via : Multiple Level Calibration



Vial: 5

Operator: GP

Inst : MS8

Multiplr: 1.00

Data File : C:\HPCHEM\1\DATA\061213J.B\J28732.D Vial: 5

Acq On : 13 Dec 06 Operator: GP 1:46 pm Sample : VBLK1213 Misc : 8260B() Inst : MS8 Multiplr: 1.00

Quant Time: Jan 3 13:27 2007

: C:\HPCHEM\1\METHODS\8260BDIR.M Method

Title : EPA METHOD 8260B

Last Update: Wed Jan 03 12:23:29 2007
Response via: Multiple Level Calibration

Response via : Multiple Level	Calibr	ation			
Internal Standards	К.Т.	QIon	Response	Conc Units	Dev(Min)
1) Fluorobenzene43) Chlorobenzene-d563) 1,4-Dichlorobenzene-d4	8.99 12.32 14.19		4363494 3446594 763801	50.00 ppb 50.00 ppb 50.00 ppb	0.00 0.00 0.01
System Monitoring Compounds 34) 1,2-Dichloroethane-d4 (S) 46) Toluene-d8 (S) 52) Bromofluorobenzene (S)	8.52 10.75 13.35	65 98 176	1249722 3894169 1317655	50.51 ppb 51.45 ppb 43.51 ppb	
Target Compounds					Ovalue

^{(#) =} qualifier out of range (m) = manual integration J28732.D 8260BDIR.M Wed Jan 03 13:27:49 2007

Spike Recovery and RPD Summary Report - WATER

: C:\HPCHEM\5\METHODS\VPHX.M

Method : C. \...
: MA VPH

Last Update : Wed Dec 20 09:14:25 2006

Response via : Initial Calibration

Non-Spiked Sample: 2121206.D

Spike

Spike

Sample

Duplicate Sample

File ID : 2121204.D

2121215.D

Sample : VLCS1212

VLCSD1212

Acq Time: 12 Dec 06 05:23 PM

13 Dec 06 10:07 AM

VPHX.M

Wed Jan 03 14:54:48 2007 GC2

FORM 4 MADEP VPH METHOD BLANK SUMMARY

VBLK1212	

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Lab File ID: 2121206.D

Lab Sample ID: VBLK1212

Matrix: (soil/water) Water Date Analyzed: 12/12/06

Instrument ID: GC2

Date Extracted:

Time Analyzed: 1845

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	LAB	CLIENT	LAB	DATE
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED
				=======================================
01	E612529-1C	841061208-27	2121208.D	12/12/06
02	E612529-2C	841061208-28	2121209.D	12/13/06
03	E612529-3C	841061208-29	2121210.D	12/13/06
04	E612529-4C	841061208-30	2121211.D	12/13/06
05	E612529-5C	841061208-31	2121212.D	12/13/06
06	E612529-6C	841061208-32	2121213.D	12/13/06
07		•	2121204.D	12/12/06
		VLCSD1212	2121215.D	12/13/06
09				
10				
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24 25				
261				
271				I
281	!			I
291				
301	l			
201		I		1

COMMENTS:	

Signal #1 : C:\HPCHEM\5\DATA\061212G2.B\2121206.D Vial: 6 Signal #2 : C:\HPCHEM\5\DATA\061212G2.B\2121206.D\2121206.D

Operator: TW

: 12 Dec 06 06:45 PM Acq On

: GC2 Inst Multiplr: 1.00

: VBLK1212 Sample

Quant Time: Dec 13 9:41 2006

: VPH()

Method

Misc

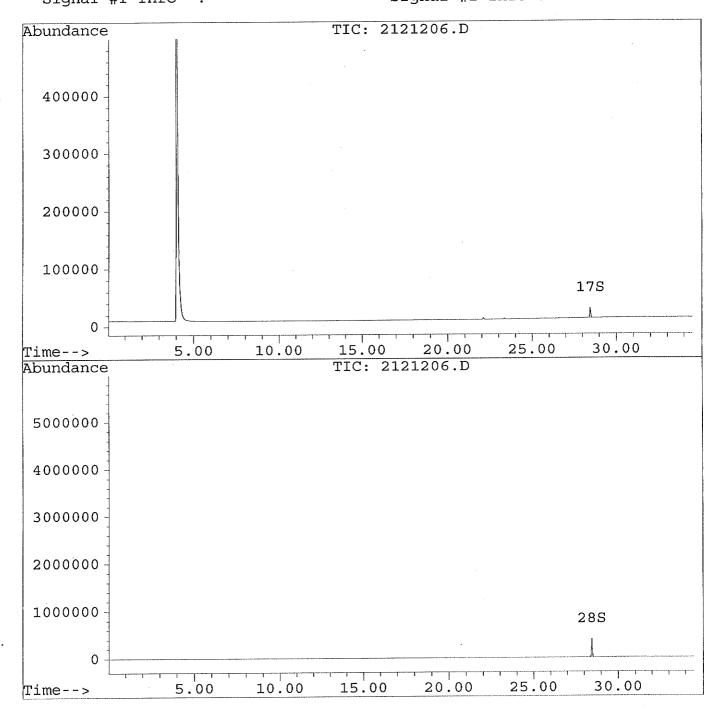
: C:\HPCHEM\5\METHODS\VPHX.M

: MA VPH Title

Last Update : Wed Dec 20 09:14:25 2006 Response via : Multiple Level Calibration

Volume Inj.

Signal #1 Phase : Signal #1 Info : Signal #2 Phase: Signal #2 Info :



GC2

Signal #1 : C:\HPCHEM\5\DATA\061212G2.B\2121206.D Vial: 6

Signal #2 : C:\HPCHEM\5\DATA\061212G2.B\2121206.D\2121206.D

Acq On : 12 Dec 06 06:45 PM Operator: TW

Inst : GC2 Sample : VBLK1212 Misc : VPH() Multiplr: 1.00

Ouant Time: Dec 13 9:41 2006

Method : C:\HPCHEM\5\METHODS\VPHX.M
Title : MA VPH

Last Update : Wed Dec 20 09:14:25 2006 Response via: Multiple Level Calibration

Volume Inj. :

Signal #2 Phase: Signal #1 Phase : Signal #2 Info: Signal #1 Info :

R.T. Response Conc Units Compound

System Monitoring Compounds System Monitoring Compounds

17) S 2,5-dibromotoluene(S) 28.47 622583 40.096 ug/L m

Recovery = 80.19%

Recovery = 80.19% 28) S 2,5-dibromotoluene(S) #2 28.46 13788612 41.642 ug/L Recovery = 83.28%

Target Compounds

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/12/06

Project No.: E612529 Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1 Location: Franklin, MA

	SPIKE	SAMPLE		QC .
	ADDED	CONCENTRATION	8	LIMITS
COMPOUND	(ug/L)	$\mid (ug/L) \mid$	REC #	REC
=====================================		=======================================	=====	=====
2-Methylnaphthalene	40.00	29.61	74	40-140
Acenaphthene	40.00	30.01	75	40-140
Acenaphthylene	40.00	31.38		40-140
Anthracene	40.00	34.82		40-140
Benzo[a]anthracene	40.00	36.49		40-140
Benzo[a]pyrene	40.00	36.35	91	40-140
Benzo[b]fluoranthene	40.00	35.67	89	40-140
Benzo[g,h,i]perylene	40.00	30.11		40-140
Benzo[k]fluoranthene	40.00	36.85		40-140
Chrysene	40.00	36.64		40-140
Decane	40.00	24.45		40-140
Dibenz[a,h]anthracene	40.00	32.14		40-140
Docosane	40.00	33.85		40-140
Dodecane	40.00	27.08		40-140
Eicosane	40.00	33.66		40-140
Fluoranthene	40.00	34.93	87	40-140
Fluorene	40.00	33.50	84	40-140
Hexacosane	40.00	34.61	86	40-140
Hexadecane	40.00	31.52		40-140
Hexatriacontane	40.00	35.66		40-140
Indeno[1,2,3-cd]pyrene	40.00	35.18	88	40-140
Naphthalene	40.00	26.75	67	40-140
Nonadecane	40.00	32.81		40-140
Nonane	40.00	25.37		30-140
Octacosane	40.00	35.82		40-140
Octadecane	40.00	31.92	80	40-140
Phenanthrene	40.00	33.62	84	40-140
Pyrene !	40.00	33.36	83	40-140
		 		

Column to be used to flag recovery values with an asterisk

COMMENTS:	

FILE: A20324.D

^{*} Values outside of QC limits

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/12/06

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1 Location: Franklin, MA

	SPIKE ADDED	SAMPLE CONCENTRATION	QC % LIMITS
COMPOUND	(ug/L)	(ug/L)	REC # REC
		=========	====== =====
Tetracosane	40.00	34.38	86 40-140
Tetradecane	40.00	29.18	73 40-140
Triacontane	40.00	36.46	91 40-140
	<u></u>		l1

COMMENTS:	

FILE: A20324.D

Page 2 of 2

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE DUPLICATE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/12/06

Project No.: E612529 Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1 Location: Franklin, MA

1	SPIKE ADDED	SAMPLE) 	0.	QC
COMPOSIND	•	· ·		8	LIMITS
COMPOUND	(ug/L)	(ug/L)	REC #		RPD REC
2-Methylnaphthalene	40.00	26.58	66	11.4	===== 25 40-140
Acenaphthene	40.00	27.18	68	9.79	25 40-140
Acenaphthylene	40.00	27.87	70	10.8	25 40-140
Anthracene	40.00	32.05	80	8.38	25 40-140
Benzo[a]anthracene	40.00	33.76	84	8.00	25 40-140
Benzo[a]pyrene	40.00	33.75	84	8.00	25 40-140
Benzo[b]fluoranthene	40.00	33.15	83	6.98	25 40-140
Benzo[g,h,i]perylene	40.00	28.60	72	4.08	25 40-140
Benzo[k]fluoranthene	40.00	34.24	86	6.74	25 40-140
Chrysene	40.00	34.08	85	7.91	25 40-140
Decane	40.00	22.31	56	8.55	25 40-140
Dibenz[a,h]anthracene	40.00	30.36	76	5.13	25 40-140
Docosane	40.00	31.55	79	7.32	25 40-140
Dodecane	40.00	24.49	61	10.8	25 40-140
Eicosane	40.00	30.99	77	8.70	25 40-140
Fluoranthene	40.00	32.11	80	8.38	25 40-140
Fluorene	40.00	30.46	76	10.0	25 40-140
Hexacosane	40.00	32.24	81	5.99	25 40-140
Hexadecane	40.00	28.64	72	9.27	25 40-140
Hexatriacontane	40.00	33.97	85	4.60	25 40-140
Indeno[1,2,3-cd]pyrene	40.00	32.89	82	7.06	25 40-140
Naphthalene	40.00	24.13	60	11.0	25 40-140
Nonadecane	40.00	30.13	75	8.92	25 40-140
Nonane	40.00	22.19	55	13.6	25 30-140
Octacosane	40.00	33.28	83	8.09	25 40-140
Octadecane	40.00	29.22	73	9.15	25 40-140
Phenanthrene	40.00	30.77	77	8.70	25 40-140
Pyrene	40.00	30.92	77	7.50	25 40-140
I					

Column to be used to flag recovery values with an asterisk

FILE: A20325.D

^{*} Values outside of QC limits

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE DUPLICATE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/12/06

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1

Location: Franklin, MA

	SPIKE	SAMPLE			QC
	ADDED	CONCENTRATION	용	%	LIMITS
COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD REC
	========	=======================================	======	======	===== =====
Tetracosane	40.00	31.91	80	7.23	25 40-140
Tetradecane	40.00	26.34	66	10.1	25 40-140
Triacontane	40.00	34.10	85	6.82	25 40-140
		<u> </u>			Iİ

 $\ensuremath{\text{\#}}$ Column to be used to flag recovery and RPD values with an asterisk

* Values outside of QC limits

COMMENTS:	·

FILE: A20325.D

Page 2 of 2

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/13/06

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1

Location: Franklin, MA

	SPIKE	SAMPLE		QC
1	ADDED	CONCENTRATION	용	LIMITS
COMPOUND	(ug/L)	(ug/L)	REC #	REC
	========		=====	======
2-Methylnaphthalene	40.00	21.35	53	40-140
Acenaphthene	40.00	22.87	57	40-140
Acenaphthylene	40.00	21.81	54	140-140
Anthracene	40.00	30.46	76	40-140
Benzo[a]anthracene	40.00	25.20	63	40-140
Benzo[a]pyrene	40.00	25.88	65	140-140
Benzo[b]fluoranthene	40.00	23.85	60	40-140
Benzo[g,h,i]perylene	40.00	24.20	60	40-140
Benzo[k]fluoranthene	40.00	26.36	66	40-140
Chrysene	40.00	23.79	59	140-140
Dibenz[a,h]anthracene	40.00	27.97	70	40-140
Fluoranthene	40.00	24.49	61	40-140
Fluorene	40.00	23.39	58	40-140
Indeno[1,2,3-cd]pyrene	40.00	22.99	57	40-140
Naphthalene	40.00	19.39	48	40-140
Phenanthrene	40.00	24.06	60	40-140
Pyrene	40.00	23.39	58	40-140

COMMENTS:		
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FILE: A20348.D

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 WATER MADEP EPH LAB CONTROL SAMPLE DUPLICATE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed: 12/13/06

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1212A-1

Location: Franklin, MA

1	L ODIKE	L CAMPIE			
	SPIKE	SAMPLE			l QC
	ADDED	CONCENTRATION	િ		LIMITS
COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD REC
======================================	========	_======================================	=====	=====	===== =====
2-Methylnaphthalene	40.00	17.58	44	18.6	25 40-140
Acenaphthene	40.00	18.71	47	19.2	25 40-140
Acenaphthylene	40.00	18.14	45	18.2	25 40-140
Anthracene	40.00	24.85	62	20.3	25 40-140
Benzo[a]anthracene	40.00	21.31	53	17.2	25 40-140
Benzo[a]pyrene	40.00	21.46	54	18.5	25 40-140
Benzo[b]fluoranthene	40.00	19.62	49	20.2	25 40-140
Benzo[g,h,i]perylene	40.00	20.31	51	16.2	25 40-140
Benzo[k]fluoranthene	40.00	22.14	55	18.2	25 40-140
Chrysene	40.00	20.21	50	16.5	25 40-140
Dibenz[a,h]anthracene	40.00	22.73	57	20.5	25 40-140
Fluoranthene	40.00	20.13	50	19.8	25 40-140
Fluorene	40.00	19.01	48	18.9	25 40-140
Indeno[1,2,3-cd]pyrene	40.00	19.02	48	17.1	25 40-140
Naphthalene	40.00	16.49	41	15.7	25 40-140
Phenanthrene	40.00	19.82	50 j	18.2	25 40-140
Pyrene	40.00	19.31 i	48	18.9	25 40-140
-			1	i	

FILE: A20349.D

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 WATER LAB CONTROL SAMPLE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed:

Project No.:

Project: 51239

Sample No.: LCS1212A-1

Location:

1		SPIKE	SAMPLE	1	QC
		ADDED	CONCENTRATION	응	LIMITS
	COMPOUND	(ug/L)	(ug/L)	REC #	REC
=		=======		======	=====
	Decane	40.00	22.25	56	40-140
	Docosane	40.00	33.21	83	40-140
	Dodecane	40.00	25.24	63	40-140
	Eicosane	40.00	35.52	89	40-140
	Hexacosane	40.00	33.33	83	40-1401
	Hexadecane	40.00	30.53	76	40-140
1	Hexatriacontane	40.00	33.03	82	40-1401
I	Nonadecane	40.00	32.01	80	40-140
	Nonane	40.00	17.87	45	30-140
	Octacosane	40.00	33.45	84	40-1401
	Octadecane	40.00	33.02	82	40-140
	Tetracosane	40.00	33.05	83 i	40-140
1	Tetradecane	40.00	28.16	70 i	40-140
	Triacontane	40.00	38.01	95 i	40-1401
١	Ì				i

COMMENTS:	

FILE: A20339.D

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 WATER LAB CONTROL SAMPLE DUPLICATE

Lab Name: PREMIER LABORATORY, LLC Date Analyzed:

Project No.:

Project: 51239

Sample No.: LCS1212A-1

Location:

ADDED CONCENTRATION % % LIMITS COMPOUND (ug/L) (ug/L) (ug/L) REC # RPD # RPD REC RPD # RPD # RPD REC RPD # RPD REC RPD # RPD REC RPD # RPD REC RPD # RPD RPD RPD REC RPD # RPD	1		ODTVO				
COMPOUND	!	· ·	SPIKE	SAMPLE		1	l QC
Decane	ı		ADDED	CONCENTRATION	용	%	LIMITS
Decane	ı	COMPOUND	(ug/L)	(ug/L)	REC #	RPD #	RPD REC
Docosane				=======================================		======	====== ======
Dodecane		Decane	40.00	19.64	49	13.3	25 40-140
Dodecane	1	Docosane	40.00	30.44	76	8.80	25 40-140
Eicosane		Dodecane	40.00	22.02	55	13.6	25 40-140
Hexacosane	1	Eicosane	40.00	32.25	81	9.41	25 40-140
Hexatriacontane		Hexacosane	40.00	30.54	76	8.80	25 40-140
Nonadecane		Hexadecane	40.00	26.98	67	12.6	25 40-140
Nonadecane		Hexatriacontane	40.00	31.05	78	5.00	25 40-140
Nonane		Nonadecane	40.00	28.56	71	11.9	25 40-140
Octacosane		Nonane	40.00	15.35	38	16.9	25 30-140
Octadecane		Octacosane	40.00	30.45	76 Ì	10.0	25 40-140
Tetracosane		Octadecane	40.00	29.96	75 j	8.92	25 40-140
	1	Tetracosane	40.00	30.13	75	10.1	25 40-140
I made a section of the section of t		Tetradecane	40.00	24.47	61 i	13.7	25 40-140
		Triacontane	40.00	31.80	80 j	17.1	25 40-140
1			ĺ	İ	ĺ		

COMMENTS:	

FILE: A20340.D

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 4 MADEP EPH METHOD BLANK SUMMARY

E	1	2	1	2	B	A	_	1	

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Lab File ID: A20323.D

Lab Sample ID: E1212BA-1

Matrix: (soil/water) Water Date Analyzed: 12/12/06

Instrument ID: GC1

Date Extracted:

Time Analyzed: 2128

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	I T A D	CLIENT	1 TAD	L Damp I
	LAB	,	LAB	DATE
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED
01	=========	ı	====================================	======
		841061208-27	A20329.D	12/13/06
	= -	1841061208-28	A20328.D	12/13/06
	•	841061208-29	A20327.D	12/12/06
			A20330.D	12/13/06
		•	A20324.D	12/12/06
	•	LCSD1212A-1	A20325.D	12/12/06
07				
08				
091				
10		[
11			i I	
12				
13				
14				
15				
16	***************************************			
17			· ·	·
181				
19		·	· 	1
20	***************************************			·
21			·	
221				
231			' '	
24				
251		<u> </u>		
26				
271	'			I
28		·		
291		<u> </u>		
301	 			l
301				

COMMENTS:		

FORM 4 MADEP EPH METHOD BLANK SUMMARY

E1212BA-1	

Project No.: E612529

Project: 20050458.B10/Nu-Style Phase II

Lab File ID: A20347.D

Lab Sample ID: E1212BA-1

Matrix: (soil/water) Water

Date Analyzed: 12/13/06

Instrument ID: GC1

Date Extracted:

Time Analyzed: 1948

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

	LAB	CLIENT	LAB	DATE
	SAMPLE NO.	SAMPLE ID	FILE ID	ANALYZED
	SAMPLE NO.		LIDE ID	ANALIZED
01	•	•	A20353.D, A2034	
	•	· ·	A20354.D, A2034	
	•		A20348.D	12/13/06
	•	LCSD1212A-1	A20349.D	12/13/06
05	•]		,,
06				
07				
80	·			
09	·			
10	·			
11				
12	` 			
13				
14 15				
16	PM	 	 	
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19	****	 		
20				
21	······································			
22				
23				
24				
25				
26				[
27				
28				
29				
301			I	

COMMENTS:	

Data File : D:\DATA\061212A.B\A20323.D

Acq On : 12 Dec 2006 9:28 pm Sample

: E1212BA-1

Operator: TW : GC1 Inst Multiplr: 1.00

Vial: 9

Misc : EPH() IntFile : AUTOINT1.E

Quant Time: Jan 3 14:58 2007 Quant Results File: TOTALTPH.RES

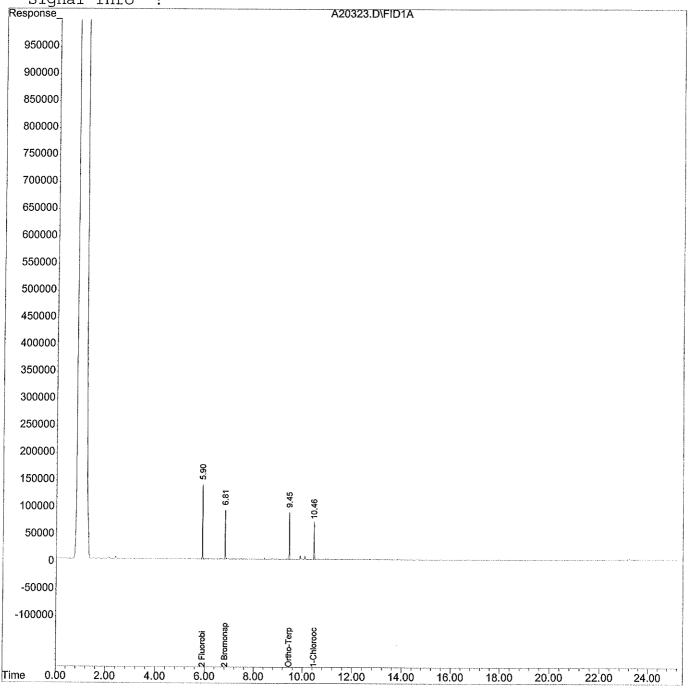
Quant Method : C:\HPCHEM\1\XMETHODS\TOTALTPH.M (Chemstation Integrator)

Title

Last Update : Wed Dec 13 12:00:38 2006 Response via: Multiple Level Calibration

DataAcq Meth : EPHACQ.M

Volume Inj. Signal Phase: Signal Info



Quantitation Report (QT Reviewed)

Vial: 9 Data File : D:\DATA\061212A.B\A20323.D Operator: TW Acq On : 12 Dec 2006 9:28 pm Acq on Sample : ETH() Inst : GC1 : E1212BA-1 Multiplr: 1.00

IntFile : AUTOINT1.E

Quant Time: Jan 3 14:58 2007 Quant Results File: TOTALTPH.RES

Quant Method : C:\HPCHEM\1\XMETHODS\TOTALTPH.M (Chemstation Integrator)

Title

Last Update : Wed Dec 13 12:00:38 2006

Response via : Initial Calibration

DataAcq Meth : EPHACQ.M

Volume Inj. : Signal Phase: Signal Info :

	Compound	l			R.	Т.	Response		Conc Units
	cem Monitor		unds			100 000 000 000 pm			
1) S	2 Fluorob	iphenyl			5.	90	1424305		36.668 ug/ml
Spiked	Amount	40.000	Range	16	-	56	Recovery	=	91.67%#
2) S	2 Bromona	phthalene	•		6.	81	1006018		38.976 ug/ml
Spiked	Amount	40.000	Range	16	-	56	Recovery	=	97.44%#
	Ortho-Ter				9.	45	1005965		23.282 ug/ml
Spiked	Amount	40.000	Range	16	_	56	Recovery	=	58.20%#
	1-Chloroc		J	:	10.	46	878640		25.818 ug/ml
Spiked	Amount	40.000	Range	16	_	56	Recovery	=	64.55%#

Target Compounds

IntFile : AUTOINT1.E

Quant Time: Dec 20 16:40 2006 Quant Results File: PAH.RES

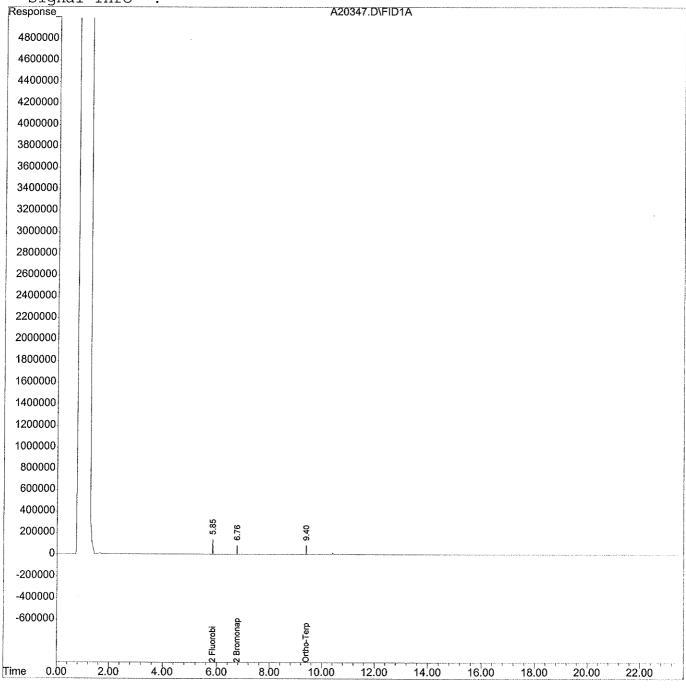
Quant Method : C:\HPCHEM\1\XMETHODS\PAH.M (Chemstation Integrator)

Title :

Last Update : Thu Dec 14 09:25:52 2006 Response via : Multiple Level Calibration

DataAcq Meth : PAHACQ.M

Volume Inj. : Signal Phase : Signal Info :



Data File : D:\DATA\061213A.B\A20347.D Vial: 9 Acq On : 13 Dec 2006 7:48 pm Operator: TW Sample : E1212BA-1 Misc : EPH() IntFile : AUTOINT1.E Inst : GC1 Multiplr: 1.00

Quant Time: Dec 20 16:40 2006 Quant Results File: PAH.RES

Quant Method : C:\HPCHEM\1\XMETHODS\PAH.M (Chemstation Integrator)

Title

Last Update : Thu Dec 14 09:25:52 2006

Response via : Initial Calibration

DataAcq Meth : PAHACQ.M

Volume Inj. : Signal Phase : Signal Info :

Compound	R.T.	Response	Conc	Units		
System Monitoring Compounds						
1) S 2 Fluorobiphenyl	5.85	1472073	30.431	ug/ml		
Spiked Amount 40.000		overy =				
2) S 2 Bromonaphthalene	6.76	940402	29.077 ug/ml			
Spiked Amount 40.000	Reco	overy =				
3) S Ortho-Terphenyl	9.40	1049620	16.251	ug/mlm		
Spiked Amount 40.000	Reco	overy =	40.63%	_		
Target Compounds						
4) TM Naphthalene	0.00	0	N.D.	ug/mld		
5) M 2-Methylnaphthalene	0.00	0	N.D.	ug/mld		
6) T Acenaphthylene	0.00	0	N.D.	ug/mld		
7) TM Acenaphthene	0.00	0	N.D.	ug/mld		
8) T Fluorene	0.00	0	N.D.	ug/mld		
9) T Phenanthrene	0.00	0	N.D.	ug/mld		
10) TM Anthracene	0.00	0	N.D.	ug/mld		
11) T Fluoranthene	0.00	0	N.D.	ug/mld		
12) TM Pyrene	0.00	0	N.D.	ug/mld		
13) T Benzo[a]anthracene	0.00	0	N.D.	ug/mld		
14) TM Chrysene	0.00	0	N.D.	ug/mld		
15) T Benzo[b] fluoranthene	0.00	0	N.D.	ug/mld		
16) T Benzo[k]fluoranthene	0.00	0	N.D.	ug/mld		
17) T Benzo[a]pyrene	0.00	0	N.D.	ug/mld		
18) T Indeno[1,2,3-cd]pyrene	0.00	0	N.D.	ug/mld		
19) T Dibenz[a,h]anthracene	0.00	0	N.D.	ug/mld		
20) T Benzo[g,h,i]perylene	0.00	0	N.D.	ug/mld		
21) x C11 - C22 Aromatics	0.00	0	N.D.	ug/mld		

IntFile : AUTOINT1.E

Quant Time: Dec 20 16:39 2006 Quant Results File: ALKEPH.RES

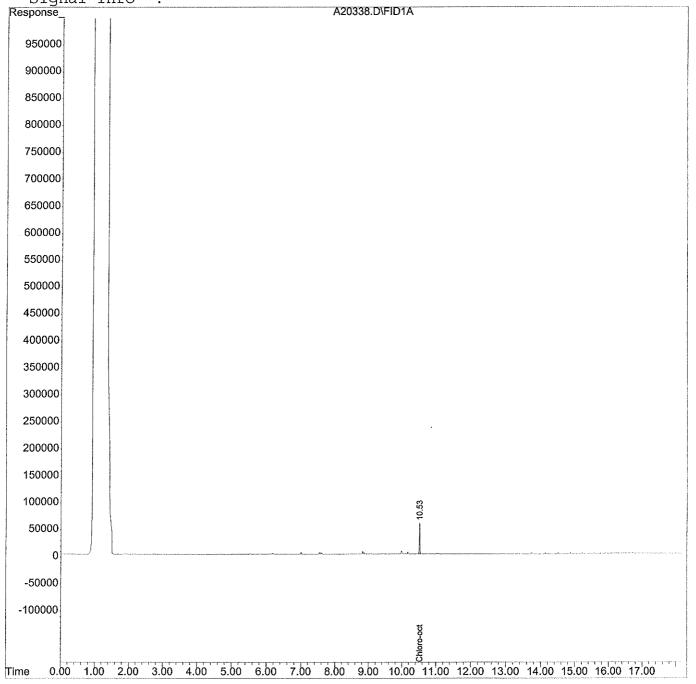
Quant Method: C:\HPCHEM\1\XMETHODS\ALKEPH.M (Chemstation Integrator)

Title :

Last Update : Thu Dec 14 09:06:29 2006 Response via : Multiple Level Calibration

DataAcq Meth : ALKNEACQ.M

Volume Inj. : Signal Phase : Signal Info :



Quantitation Report (QT Reviewed)

Data File : D:\DATA\061213A.B\A20338.D

Acq On : 13 Dec 2006 4:02 pm Sample : E1212BA-1 Misc : EPH() IntFile : AUTOINT1.E

Vial: 2 Operator: TW

Inst : GC1 Multiplr: 1.00

Ouant Time: Dec 20 16:39 2006 Ouant Results File: ALKEPH.RES

Quant Method: C:\HPCHEM\1\XMETHODS\ALKEPH.M (Chemstation Integrator) Title

Last Update : Thu Dec 14 09:06:29 2006

Response via: Initial Calibration

DataAcq Meth : ALKNEACQ.M

Volume Inj. : Signal Phase : Signal Info :

Compound

R.T. Response Conc Units

1) S Chloro-octadecane * 10.53 649281 18.615 ug/ml piked Amount 40.000 Recovery - 10.53

Spiked Amount 40.000

Target Compounds

ICP Method Blank Summary

Workorder#:	E612529	Matrix: Aque	
Element	Result	MDL	Run Date
	ug/L	ug/L	
Antimony	ND	10	12/18/2006
Aluminum		50	Name of the last o
Arsenic	ND	10	12/18/2006
Barium	ND	10	
Beryllium	ND	1	12/18/2006
Boron		10	
Calcium		50	
Cadmium	ND	2	12/18/2006
Cobalt		2	
Chromium	ND	10	
Copper	ND	10	12/18/2006
Iron		50	
Lead	ND	4	12/18/2006
Magnesium		50	
Manganese		10	
Mercury CV	ND	0.2	12/14/2006
Molybdenum		10	
Nickel	ND	10	12/18/2006
Potassium		2000	Commence of American Specials and Commence of Commence
Selenium	ND	10	
Silver	ND	2	12/18/2006
Sodium		1000	
Thallium	ND	5	12/18/2006
Titanium		2	
Tin		50	
Vanadium		10	
Zinc	ND	10	12/18/2006

ND = NONE DETECTABLE

- (*) Elevated MDLs due to dilution for range
- (**) Elevated MDLs due to dilution for interferences

Fortified Sample/Blank Recovery Report

Date:	December 18, 2006
Time:	9:19
Method:	200.7/6010B
Analyst:	BSZ
Validator	•

	Matrix	•
Aqueous		
TCLP		
SPLP		

Run Log Reference Number TRC1218-1

Workorder #: E612529

Fortified Sample ID #:

E61252711C

Units: ug/L

				Matrix(LFM)/LFM Duplicate								LFB	
Element	Sample	Sample	RPD	Spike	LFM	%	LFMD	%	Recovery	RPD	Result	%	Recovery
		Duplicate		Amount	Result	recovery	Result	recovery	Limits			recovery	Limits
Ag	0	0		500	507.9	101.6	510	102.0	75-125	0.4	510.2	102.0	80-120
Al				10000					75-125		***		80-120
As	0	0		500	486.7	97.3	483.2	96.6	75-125	0.7	485	97.0	80-120
В				500					75-125				80-120
Ва	56.98	57	0.0	500	574.9	103.6	574.4	103.5	75-125	0.1	526.2	105.2	80-120
Ве	0	0		500	500	100.0	501.4	100.3	75-125	0.3	514.4	102.9	
Ca				11000					75-125			****	80-120
Cd	0	0		500	501.9	100.4	504.6	100.9	75-125	0.5	518	103.6	80-120
Co				500					75-125				80-120
Cr	0	0		500	492.4	98.5	495	99.0	75-125	0.5	506.5	101.3	80-120
Cu	0	0		500	495.9	99.2	497.2	99.4	75-125	0.3	502.1	100.4	80-120
Fe				500				******	75-125	***************************************		****	80-120
<				25000			***		75-125				80-120
Иg				11000				***************************************	75-125			****	80-120
Mn				500					75-125		****		80-120
VIO				500		*****			75-125		***************************************		80-120
Na				11000		Ì			75-125				80-120
Vi	0	0		500	501.4	100.3	501.3	100.3	75-125	0.0	513.1	102.6	80-120
⊃b	0	0		500	477.4	95.5	482.7	96.5	75-125	1.1	493.7	98.7	80-120
Sb	0	0		500	464.4	92.9	457.9	91.6	75-125	1.4	483.2	96.6	80-120
Se	0	0		500	486.5	97.3	491.9	98.4	75-125	1.1	497.3	99.5	80-120
Sn Sn	Ì			2500					75-125			55.0	80-120
П	0	0		500	498.7	99.7	496.1	99.2	75-125	0.5	513	102.6	80-120
Гi				500				· -	75-125				80-120
/				500					75-125				80-120
Zn	32.76	33.58	2.5	500	504.4	94.3	508.2	95.1	75-125	0.8	492.6	98.5	80-120
∃g	0	0		5	4.839	96.8	4.861	97.2	66-133	0.5	4.955	99.1	80-120

COMMENTS: saplme spiked for Hg: E61257801



Modified Tier I Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E612052
Date Samples Received at Laboratory:	12/1/2006
Date of Review:	1/11/2007

Seventeen soil samples, including one field duplicate, were collected and submitted to Premier Laboratory, LLC in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant metals plus barium by EPA Methods 6010B and 7471, cyanide by EPA Method 9012, polychlorinated biphenyls (PCBs) by EPA Method 8082, and petroleum hydrocarbons by Massachusetts Department of Environmental Protection (MADEP) Methods Extractable Petroleum Hydrocarbons (EPH) and Volatile Petroleum Hydrocarbons (VPH). One aqueous trip blank was also submitted for analysis of VOCs by EPA Method 8260B. Dedicated sampling equipment was employed; therefore, no equipment blank was indicated.

Samples were analyzed within method-specified holding times and in accordance with the Massachusetts Contingency Plan (MCP) Compendium of Analytical Methods (CAM) data enhancement protocols.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

4 043 575 57 50 45 57 57 57 57 57 57 57 57 57 57 57 57 57		<u>YES</u>	N	O
1. SAMPLING AND FIELD MEASUREM	ENTS:			
Field measurement calibration records				
Groundwater field measurements (if applicable)				O NA
Soil sampling field measurements (if applicable)		₩ W		
Sediment sampling field measurements (if applicable)			;	o NA
Surface water sampling field measurements (if applica	ıble)			ONA
Low-flow sampling field measurements (if applicable)	1			DNA
Documentation of field activities		A R R R C		
Sample numbering and labeling		区		
Chain-of-Custody records		図		
Trip blanks		. 🗹 .	1	
Duplicate samples		□		
Equipment blanks			Ī]NA
Split samples (if any)				JNA
2. LABORATORY MEASUREMENTS:		,		
Trip blanks		⊡′,	[
Instrument blanks				
Laboratory control samples		Ū	[
Duplicates samples		년 년	[Ī
Equipment blanks]] ~A
Matrix spike/matrix spike duplicates		Ø .		
Analysis type				<u> </u>
Chain-of-Custody records			Γ	
Surrogate recoveries			Ē	<u>-</u>
Sample Project Narratives		- FT	Ē	Ī
Split samples (if any)	•			J N/A
		1.1	6	- /:
	TOTAL:	16		-
	PERCENT C	OMDI ETE	100	0./
	ETTUCEINI (OMPLETE:	, , - '	%



Premier Laboratory, LLC

61 Louisa Viens Drive Dayville, CT 06241

Telephone: 860-774-6814 Fax: 860-774-26A9

ANALYTICAL DATA & QUALITY CONTROL REPORT

Report Number: E612052

Project: 20050458.B10/Nu-Style Phase II

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss





ANALYTICAL DATA REPORT

Report Number: E612052 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

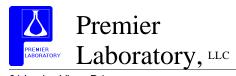
Received Date: 12/1/2006 Report Date: 3/13/2007

Premier Laboratory, LLC Authorized Signature



956D450

Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)



	MADEP MCP Analytical Method Report Certification Form									
Labo	ratory Name: Pre	emier Laboratory, L	LC		Project #:	E612	052			
	ct Location: Fran				MADEP R					
			llowing data set:[list	Laboratory S	ample ID Numb	er(s)]				
1, 10	, 11, 12, 13, 14, 1	5, 16, 17, 18, 2, 3,	4, 5, 6, 7, 8, 9							
	Sample Matrices: Groundwater Soil/Sediment Drinking Water Other									
	SW-846	8260B ⊠	8151A □	833		10B ⊠			0A/1A ⊠	
Meth	ods Used	8270C 🗆	8081A □			020 🗆			14M ² 🗆	
	ecified in MADEP	8082 🗵	8021B 🗆		H⊠ 7000	S ³ □		7	7196A □	
	endium of ical Methods.		ing Number (RTN), if kn d 9014 or MADEP Physi		ble Cvanide (PAC)	Method				
	all that apply)		ds 7000 Series List indi							
	An affirmative re	esponse to questic	ons A, B, C, and D	is required fo	or "Presumptiv	e Certa	inty'	' sta	tus	
Α	Were all samples	s received by the la	boratory in a condit	ion consisten	t with		Yes		No ¹	
	that described or	the Chain-of-Cust	ody documentation	for the data s	et?					
В	Were all QA/QC	procedures require	d for the specified a	nalytical meth	hod(s)	\boxtimes	Yes		No ¹	
	included in this re	eport followed, inclu	iding the requiremer	nt to note and	I					
	discuss in a narra	ative QC data that	did not meet approp	riate performa	ance					
	standards or guid	delines?								
С	Does the analytic	cal data included in	this report meet all	the requirem	ents		Yes		No ¹	
	for "Presumptive	Certainty", as desc	cribed in Section 2.0	(a),(b),(c) an	d (d) of the					
	MADEP docume	nt CAM VII A, "Qua	ality Assurance and	Quality Contr	rol Guidelines					
	for the Acquisitio	n and Reporting of	Analytical Data"?							
D	VPH and EPH M	lethods only: Was	s the VPH or EPH m	ethod run wit	thout		Yes		No ¹	
	significant modifi	cations, as specifie	ed in Section 11.3?							
	A respons	e to questions E a	and F below is requ	ired for "Pre	sumptive Cert	ainty" s	statu	s		
E	Were all QC perf	ormance standards	and recommendation	ons for the			Yes		No ¹	
	specified method	s achieved?								
F	Were results for	all analyte-list com	pounds/elements for	the specified	d		Yes		No ¹	
	method(s) reporte	ed?								
	¹ All NO ans	wers must be addre	essed in an attached	d Environmen	tal Laboratory o	case na	rrativ	e.		
I tho	undereigned ettect u	nder the naine and not	nalties of perjury that, ba	and upon my n	organal					
			rmation, the material co							
analy	tical report is, to the b	est of my knowledge a	and belief, accurate and	l complete.						
Sign	ature:	William.		Position:	Laboratory D	irector				
	= .			.						
Print	Printed Name: Robert Stevenson Date: 3/13/2007									

Page 2 of 103

> Report No: E612052 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory received 18 samples from Fuss & O'Neill on 12/01/2006. The samples were analyzed from the following list of analyses:

Cyanide, Total, by 9012 in GW/SW 9012[9012]
Moisture, Percent
Trace Priority Pollutant (13) Metals in Soil 6010B[3000], 7471[7471]
Volatiles by 8260B in GW/SW 8260B

Extractable Petroleum Hydrocarbon (EPH)
MADEP EPH[MADEP EPH]
PCB's by 8082 in GW/SW
8082[3500]
Volatile Petroleum Hydrocarbon (VPH)
MADEP VPH

In order to meet requested detection limits, EDB results were estimated to 3 ppb for EPA method 8260B. Dibromochloromethane, 1,2-Dichlorobenzene and 1,1,2,2-Tetrachloroethane were all estimated to a value of 5.0 ppb. This value of 5.0 ppb corresponds to the lowest level of calibration on the instrument prior to the % solid value being calculated into the reported detection limits. The samples were ND for all estimated compounds to their respective values.

Variances:

SDG:

A full list 8260B LCS was run and met the applicable recovery criteria for "Presumptive Certainty". An LCS Duplicate encompasing all target compounds was not run for EPA method 8260B. Both an LCS and LCSD were analyzed for the oxygenate compounds only.

Method:

None reported.

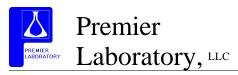
QA/QC:

956D450

Sample 12A, 841061130-12, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 13A, 841061130-13, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

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> Report No: E612052 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY (continued)

QA/QC (continued):

Sample 13A, 841061130-13, Volatiles by 8260B: Two surrogate spikes were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

Sample 14A, 841061130-14, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 15A, 841061130-15, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 15A, 841061130-15, Volatiles by 8260B: Two surrogate spikes were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

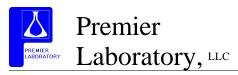
Sample 16A, 841061130-16, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 16A, 841061130-16, Volatiles by 8260B: Two surrogate spikes were outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

Sample 17A, 841061130-17, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 1A, 841061130-01, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

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61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

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Report No: E612052 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY (continued)

QA/QC (continued):

Sample 3A, 841061130-03, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 5A, 841061130-05, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 5A, 841061130-05, Volatiles by 8260B: One surrogate spike was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the surrogate was still outside of the limits.

Sample 6A, 841061130-06, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

Sample 8A, 841061130-08, Volatiles by 8260B: The matrix spike/ matrix spike duplicate recoveries for the sample were outside of the established control limits due to matrix interference. The associated LCS recoveries were within the established quality control limits.

Sample 9A, 841061130-09, Volatiles by 8260B: One internal standard was outside quality control limits for the sample due to matrix interference. The sample was re-analyzed and the internal standard was still outside the limits.

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(1) 0410/1130 01					
(1) 841061130-01	-4 C-11-J				
	atrix: Solid	0.52	/1	12/06/06 1	2.11 DDD
Cyanide, Total, by SW-846 9012	ND	0.53	mg/kg	12/06/06 1	2:11 DDD
Trace Metals by 6010B	ND	0.52	Л	12/05/06	DCZ
Antimony	ND	0.53	mg/kg	12/05/06	BSZ
Arsenic	ND	0.50	mg/kg	12/05/06	BSZ
Barium	16	0.50	mg/kg	12/05/06	BSZ
Beryllium	0.19	0.050	mg/kg	12/05/06	BSZ
Cadmium	ND	0.10	mg/kg	12/05/06	BSZ
Chromium	3.2	0.50	mg/kg	12/05/06	BSZ
Copper	4.9	0.50	mg/kg	12/05/06	BSZ
Lead	4.9	0.20	mg/kg	12/05/06	BSZ
Nickel	2.6	0.50	mg/kg	12/05/06	BSZ
Selenium	ND	0.50	mg/kg	12/05/06	BSZ
Silver	ND	0.10	mg/kg	12/05/06	BSZ
Thallium	ND	0.26	mg/kg	12/08/06	BSZ
Zinc	10	0.50	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.021	mg/kg	12/05/06	AM
(2) 841061130-02					
Date Collected: 11/30/2006 M	atrix: Solid				
Cyanide, Total, by SW-846 9012	ND	0.56	mg/kg	12/06/06 1	2:12 DDD
Trace Metals by 6010B					
Antimony	ND	0.56	mg/kg	12/05/06	BSZ
Arsenic	ND	0.56	mg/kg	12/05/06	BSZ
Barium	24	0.56	mg/kg	12/05/06	BSZ
Beryllium	0.57	0.056	mg/kg	12/05/06	BSZ
Cadmium	ND	0.11	mg/kg	12/05/06	BSZ
Chromium	5.3	0.56	mg/kg	12/05/06	BSZ
Copper	12	0.56	mg/kg	12/05/06	BSZ
Lead	8.1	0.22	mg/kg	12/05/06	BSZ
Nickel	3.3	0.56	mg/kg	12/05/06	BSZ
Selenium	ND	0.56	mg/kg	12/05/06	BSZ
Silver	ND	0.11	mg/kg	12/05/06	BSZ
Thallium	ND	0.28	mg/kg	12/08/06	BSZ
Zinc	13	0.56	mg/kg	12/05/06	BSZ

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(2) 841061130-02 (continued) Date Collected: 11/30/2006 Matrix: Solid					
Mercury by SW-846 7471 in SW	ND	0.022	mg/kg	12/05/06	AM
(3) 841061130-03					
Date Collected: 11/30/2006 Matrix: Solid					
Cyanide, Total, by SW-846 9012	ND	0.56	mg/kg	12/06/06 12:13	DDD
Trace Metals by 6010B					
Antimony	ND	0.56	mg/kg	12/05/06	BSZ
Arsenic	1.2	0.56	mg/kg	12/05/06	BSZ
Barium	36	0.56	mg/kg	12/05/06	BSZ
Beryllium	0.21	0.056	mg/kg	12/05/06	BSZ
Cadmium	0.17	0.11	mg/kg	12/05/06	BSZ
Chromium	7.1	0.56	mg/kg	12/05/06	BSZ
Copper	91	0.56	mg/kg	12/05/06	BSZ
Lead	40	0.22	mg/kg	12/05/06	BSZ
Nickel	4.0	0.56	mg/kg	12/05/06	BSZ
Selenium	ND	0.56	mg/kg	12/05/06	BSZ
Silver	ND	0.11	mg/kg	12/05/06	BSZ
Thallium	ND	0.28	mg/kg	12/05/06	BSZ
Zinc	85	0.56	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	0.029	0.022	mg/kg	12/05/06	AM
(4) 841061130-04					
Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012	ND	0.59	mg/kg	12/06/06 12:14	DDD

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(4) 841061130-04 (continued)					
Date Collected: 11/30/2006 Matrix: Solid	l				
Trace Metals by 6010B					
Antimony	ND	0.59	mg/kg	12/05/06	BSZ
Arsenic	ND	0.59	mg/kg	12/05/06	BSZ
Barium	20	0.59	mg/kg	12/05/06	BSZ
Beryllium	0.19	0.059	mg/kg	12/05/06	BSZ
Cadmium	0.14	0.12	mg/kg	12/05/06	BSZ
Chromium	6.0	0.59	mg/kg	12/05/06	BSZ
Copper	43	0.59	mg/kg	12/05/06	BSZ
Lead	18	0.24	mg/kg	12/05/06	BSZ
Nickel	3.6	0.59	mg/kg	12/05/06	BSZ
Selenium	ND	0.59	mg/kg	12/05/06	BSZ
Silver	ND	0.12	mg/kg	12/05/06	BSZ
Thallium	ND	0.30	mg/kg	12/05/06	BSZ
Zinc	63	0.59	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.024	mg/kg	12/05/06	AM
(5) 841061130-05					
Date Collected: 11/30/2006 Matrix: Solid	[
Cyanide, Total, by SW-846 9012	ND	0.59	mg/kg	12/06/06	12:15 DDD
Trace Metals by 6010B					
Antimony	ND	0.59	mg/kg	12/05/06	BSZ
Arsenic	6.6	0.59	mg/kg	12/05/06	BSZ
Barium	36	0.59	mg/kg	12/05/06	BSZ
Beryllium	0.22	0.059	mg/kg	12/05/06	BSZ
Cadmium	0.13	0.12	mg/kg	12/05/06	BSZ
Chromium	35	0.59	mg/kg	12/05/06	BSZ
Copper	160	0.59	mg/kg	12/05/06	BSZ
Lead	25	0.23	mg/kg	12/05/06	BSZ
Nickel	6.2	0.59	mg/kg	12/05/06	BSZ
Selenium	ND	0.59	mg/kg	12/05/06	BSZ
Silver	ND	0.12	mg/kg	12/05/06	BSZ
Thallium	ND	0.29	mg/kg	12/08/06	BSZ
Zinc	27	0.59	mg/kg	12/05/06	BSZ
Mercury by SW-846 7471 in SW	0.14	0.023	mg/kg	12/05/06	AM

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Company	Parameter	Result	DL	Units	Completed	By Dilution
National National	(6) 9410(1120.0)					
Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 1206/06 12:16 DDD Trace Metals by 6010B ND 0.57 mg/kg 1205/06 BSZ Arsenic 2.6 0.57 mg/kg 1205/06 BSZ Barium 36 0.57 mg/kg 1205/06 BSZ Beryllium 0.15 0.057 mg/kg 1205/06 BSZ Cadmium ND 0.11 mg/kg 1205/06 BSZ Cadmium ND 0.11 mg/kg 1205/06 BSZ Chromium 4.1 0.57 mg/kg 1205/06 BSZ Copper 9.0 0.57 mg/kg 1205/06 BSZ Lead 89 0.23 mg/kg 1205/06 BSZ Selenium ND 0.57 mg/kg 1205/06 BSZ Silver ND 0.11 mg/kg 1205/06 BSZ Thallium ND 0.23 mg/kg 1205/06<		wire Colid				
Trace Metals by 6010B			0.57	ma/ka	12/06/06 12:16	מממ
Antimony Antimony Antimony Arsenic 2.6 Arsenic 2.6 0.57 Argkg 1205/06 BSZ Barium 3.6 0.57 Barium 3.6 0.57 Barium 3.6 0.57 BSZ Beryllium 0.15 0.057 BSZ Beryllium 0.15 0.057 BSZ Beryllium ND 0.11 Bykg 1205/06 BSZ Cadmium ND 0.11 Bykg 1205/06 BSZ Cadmium ND 0.11 Bykg 1205/06 BSZ Commium 4.1 0.57 Bykg 1205/06 BSZ Commium 4.1 0.57 Bykg 1205/06 BSZ Copper 9.0 0.57 Bykg 1205/06 BSZ Copper 9.0 0.57 Bykg 1205/06 BSZ Copper 10.0 BSZ Copper 9.0 0.57 Bykg 1205/06 BSZ Copper ND 0.57 Bykg 1205/06 BSZ Copper ND 0.57 Bykg 1205/06 BSZ Copper ND 0.11 Bykg 1205/06 BSZ Copper ND 0.11 Bykg 1205/06 BSZ Copper ND 0.11 Bykg 1205/06 BSZ Copper ND 0.11 Bykg 1205/06 BSZ Copper ND 0.28 Bykg 1205/06 BSZ Copper Date Collected: 11/30/2006 Matrix: Solid ND 0.57 Bykg 1205/06 BSZ Arsenic ND 0.57 Bykg 1205/06 BSZ Copper ND 0.57 Bykg 1205/06 BSZ Copper 13 0.57 Bykg 1208/06 BSZ Copper 13	•	ND	0.57	mg/kg	12/00/00 12.10	עטט
Arsenic 2.6	•	MD	0.57	ma/ka	12/05/06	RC7
Barium	•					
Beryllium						
Cadmium ND 0.11 mg/kg 12/05/06 BSZ Chromium 4.1 0.57 mg/kg 12/05/06 BSZ Copper 9.0 0.57 mg/kg 12/05/06 BSZ Lead 89 0.23 mg/kg 12/05/06 BSZ Nickel 5.0 0.57 mg/kg 12/05/06 BSZ Selenium ND 0.57 mg/kg 12/05/06 BSZ Silver ND 0.11 mg/kg 12/05/06 BSZ Thallium ND 0.28 mg/kg 12/05/06 BSZ Zinc 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 BSZ Mercury by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B Matrix: Solid Solid Solid BSZ Arsenic ND 0.57 mg/kg 12/0						
Chromium						
Copper 9.0 0.57 mg/kg 12/05/06 BSZ Lead 89 0.23 mg/kg 12/05/06 BSZ Nickel 5.0 0.57 mg/kg 12/05/06 BSZ Selenium ND 0.57 mg/kg 12/05/06 BSZ Silver ND 0.11 mg/kg 12/05/06 BSZ Thallium ND 0.28 mg/kg 12/05/06 BSZ Zinc 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 BSZ Mercury by SW-846 9012 ND 0.57 mg/kg 12/05/06 AM Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B Antimony ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ						
Lead 89						
Nickel						
Selenium ND 0.57 mg/kg 12/05/06 BSZ Silver ND 0.11 mg/kg 12/05/06 BSZ Thallium ND 0.28 mg/kg 12/05/06 BSZ Zine 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 AM (7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B Antimony ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ						
Silver ND 0.11 mg/kg 12/05/06 BSZ Thallium ND 0.28 mg/kg 12/05/06 BSZ Zinc 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 AM (7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B Antimony ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ						
Thallium ND 0.28 mg/kg 12/05/06 BSZ Zinc 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 AM (7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.						
Zinc 54 0.57 mg/kg 12/05/06 BSZ Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 AM (7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 <						
Mercury by SW-846 7471 in SW ND 0.023 mg/kg 12/05/06 AM (7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND						
(7) 841061130-07 Date Collected: 11/30/2006 Matrix: Solid						
Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ	Welculy by 5 W-040 7471 III 5 W	ND	0.023	mg/kg	12/03/00	AWI
Date Collected: 11/30/2006 Matrix: Solid Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ	(7) 841061130-07					
Cyanide, Total, by SW-846 9012 ND 0.57 mg/kg 12/06/06 12:18 DDD Trace Metals by 6010B ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg		rix: Solid				
Trace Metals by 6010B Antimony ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ			0.57	mg/kg	12/06/06 12:18	DDD
Antimony ND 0.57 mg/kg 12/08/06 BSZ Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	•			88		
Arsenic ND 0.57 mg/kg 12/08/06 BSZ Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	•	ND	0.57	mg/kg	12/08/06	BSZ
Barium 48 0.57 mg/kg 12/08/06 BSZ Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	•					
Beryllium 0.34 0.057 mg/kg 12/08/06 BSZ Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Barium					
Cadmium 0.34 0.11 mg/kg 12/08/06 BSZ Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Beryllium	0.34			12/08/06	BSZ
Chromium 8.4 0.57 mg/kg 12/08/06 BSZ Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	•	0.34			12/08/06	BSZ
Copper 13 0.57 mg/kg 12/08/06 BSZ Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Chromium	8.4	0.57		12/08/06	BSZ
Lead 8.4 0.23 mg/kg 12/08/06 BSZ Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Copper	13	0.57		12/08/06	BSZ
Nickel 23 0.57 mg/kg 12/08/06 BSZ Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	* *	8.4	0.23		12/08/06	BSZ
Selenium ND 0.57 mg/kg 12/08/06 BSZ Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Nickel	23			12/08/06	BSZ
Silver ND 0.11 mg/kg 12/08/06 BSZ Thallium ND 0.28 mg/kg 12/08/06 BSZ	Selenium					
Thallium ND 0.28 mg/kg 12/08/06 BSZ	Silver					
	Thallium	ND	0.28		12/08/06	BSZ
Zinc 20 0.57 mg/kg 12/08/00 BSZ	Zinc	20	0.57	mg/kg	12/08/06	BSZ

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(7) 841061130-07 (continued) Date Collected: 11/30/2006 Matrix:				40.00	
Mercury by SW-846 7471 in SW	0.034	0.023	mg/kg	12/05/06	AM
(8) 841061130-08					
Date Collected: 11/30/2006 Matrix:	<u>Solid</u>				
Cyanide, Total, by SW-846 9012	ND	0.57	mg/kg	12/06/06 12:19	DDD
Trace Metals by 6010B					
Antimony	ND	0.57	mg/kg	12/08/06	BSZ
Arsenic	2.0	0.57	mg/kg	12/08/06	BSZ
Barium	24	0.57	mg/kg	12/08/06	BSZ
Beryllium	0.36	0.057	mg/kg	12/08/06	BSZ
Cadmium	0.19	0.11	mg/kg	12/08/06	BSZ
Chromium	5.4	0.57	mg/kg	12/08/06	BSZ
Copper	18	0.57	mg/kg	12/08/06	BSZ
Lead	22	0.23	mg/kg	12/08/06	BSZ
Nickel	37	0.57	mg/kg	12/08/06	BSZ
Selenium	ND	0.57	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.28	mg/kg	12/08/06	BSZ
Zinc	26	0.57	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	0.051	0.023	mg/kg	12/05/06	AM
(9) 841061130-09					
Date Collected: 11/30/2006 Matrix: Cyanide, Total, by SW-846 9012	<mark>Solid</mark> ND	0.53	mg/kg	12/06/06 12:20	DDD

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(9) 841061130-09 (continued)					
Date Collected: 11/30/2006 Matrix: Solid					
Trace Metals by 6010B					
Antimony	ND	0.53	mg/kg	12/08/06	BSZ
Arsenic	1.1	0.53	mg/kg	12/08/06	BSZ
Barium	39	0.53	mg/kg	12/08/06	BSZ
Beryllium	0.25	0.053	mg/kg	12/08/06	BSZ
Cadmium	0.22	0.10	mg/kg	12/08/06	BSZ
Chromium	5.1	0.53	mg/kg	12/08/06	BSZ
Copper	32	0.53	mg/kg	12/08/06	BSZ
Lead	20	0.21	mg/kg	12/08/06	BSZ
Nickel	4.9	0.53	mg/kg	12/08/06	BSZ
Selenium	ND	0.53	mg/kg	12/08/06	BSZ
Silver	ND	0.10	mg/kg	12/08/06	BSZ
Thallium	ND	0.26	mg/kg	12/08/06	BSZ
Zinc	48	0.53	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	0.023	0.021	mg/kg	12/05/06	AM
(10) 841061130-10					
Date Collected: 11/30/2006 Matrix: Solid					
Cyanide, Total, by SW-846 9012	ND	0.54	mg/kg	12/06/06	12:21 DDD
Trace Metals by 6010B					
Antimony	ND	0.54	mg/kg	12/08/06	BSZ
Arsenic	ND	0.54	mg/kg	12/08/06	BSZ
Barium	10	0.54	mg/kg	12/08/06	BSZ
Beryllium	0.13	0.054	mg/kg	12/08/06	BSZ
Cadmium	ND	0.11	mg/kg	12/08/06	BSZ
Chromium	3.5	0.54	mg/kg	12/08/06	BSZ
Copper	3.4	0.54	mg/kg	12/08/06	BSZ
Lead	1.6	0.22	mg/kg	12/08/06	BSZ
Nickel	14	0.54	mg/kg	12/08/06	BSZ
Selenium	ND	0.54	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.27	mg/kg	12/08/06	BSZ
Zinc	6.8	0.54	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.022	mg/kg	12/05/06	AM

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(11) 841061130-11					
Date Collected: 11/30/2006 Matrix: Solid	1				
Cyanide, Total, by SW-846 9012	ND	0.54	mg/kg	12/06/06 12:22	DDD
Trace Metals by 6010B	1,2	o.e .		12/00/00 12:22	222
Antimony	ND	0.54	mg/kg	12/11/06	BSZ
Arsenic	ND	0.54	mg/kg	12/08/06	BSZ
Barium	28	0.54	mg/kg	12/08/06	BSZ
Beryllium	0.18	0.054	mg/kg	12/08/06	BSZ
Cadmium	0.46	0.11	mg/kg	12/08/06	BSZ
Chromium	5.8	0.54	mg/kg	12/08/06	BSZ
Copper	31	0.54	mg/kg	12/08/06	BSZ
Lead	97	0.21	mg/kg	12/08/06	BSZ
Nickel	10	0.54	mg/kg	12/08/06	BSZ
Selenium	ND	0.54	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.27	mg/kg	12/08/06	BSZ
Zinc	71	0.54	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.021	mg/kg	12/05/06	AM
(12) 841061130-12					
Date Collected: 11/30/2006 Matrix: Solid	ļ				
Cyanide, Total, by SW-846 9012	ND	0.59	mg/kg	12/06/06 12:23	DDD
Trace Metals by 6010B					
Antimony	ND	0.59	mg/kg	12/08/06	BSZ
Arsenic	1.8	0.59	mg/kg	12/11/06	BSZ
Barium	26	0.59	mg/kg	12/08/06	BSZ
Beryllium	0.16	0.059	mg/kg	12/08/06	BSZ
Cadmium	0.13	0.12	mg/kg	12/08/06	BSZ
Chromium	7.4	0.59	mg/kg	12/08/06	BSZ
Copper	20	0.59	mg/kg	12/08/06	BSZ
Lead	25	0.23	mg/kg	12/11/06	BSZ
Nickel	2.6	0.59	mg/kg	12/08/06	BSZ
Selenium	ND	0.59	mg/kg	12/08/06	BSZ
Silver	ND	0.12	mg/kg	12/08/06	BSZ
Thallium	ND	0.29	mg/kg	12/08/06	BSZ
Zinc	14	0.59	mg/kg	12/08/06	BSZ

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(12) 841061130-12 (continued) Date Collected: 11/30/2006 Matrix: Sol Mercury by SW-846 7471 in SW	id 0.065	0.023	mg/kg	12/05/06	AM
(13) 841061130-13 Date Collected: 11/30/2006 Matrix: Sol Cyanide, Total, by SW-846 9012 Trace Metals by 6010B	<u>id</u> ND	0.55	mg/kg	12/06/06	DDD
Antimony	ND	0.55	mg/kg	12/08/06	BSZ
Arsenic	ND	0.55	mg/kg	12/08/06	BSZ
Barium	18	0.55	mg/kg	12/08/06	BSZ
Beryllium	0.12	0.055	mg/kg	12/08/06	BSZ
Cadmium	0.16	0.11	mg/kg	12/08/06	BSZ
Chromium	2.2	0.55	mg/kg	12/08/06	BSZ
Copper	5.0	0.55	mg/kg	12/08/06	BSZ
Lead	9.2	0.22	mg/kg	12/11/06	BSZ
Nickel	3.2	0.55	mg/kg	12/08/06	BSZ
Selenium	ND	0.55	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.27	mg/kg	12/08/06	BSZ
Zinc	14	0.55	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.022	mg/kg	12/05/06	AM
(14) 841061130-14 <u>Date Collected: 11/30/2006</u> Cyanide, Total, by SW-846 9012 Matrix: Sol	<u>id</u> ND	0.55	mg/kg	12/06/06	DDD

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(14) 841061130-14 (continued)					
· · ·	atrix: Solid				
Trace Metals by 6010B					
Antimony	ND	0.55	mg/kg	12/08/06	BSZ
Arsenic	ND	0.55	mg/kg	12/08/06	BSZ
Barium	11	0.55	mg/kg	12/08/06	BSZ
Beryllium	0.17	0.055	mg/kg	12/08/06	BSZ
Cadmium	ND	0.11	mg/kg	12/11/06	BSZ
Chromium	5.5	0.55	mg/kg	12/08/06	BSZ
Copper	2.9	0.55	mg/kg	12/08/06	BSZ
Lead	2.6	0.22	mg/kg	12/11/06	BSZ
Nickel	1.8	0.55	mg/kg	12/08/06	BSZ
Selenium	ND	0.55	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.27	mg/kg	12/08/06	BSZ
Zinc	6.3	0.55	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.022	mg/kg	12/05/06	AM
(15) 841061130-15					
` '	atrix: Solid				
Cyanide, Total, by SW-846 9012	ND	0.56	mg/kg	12/06/06	DDD
Trace Metals by 6010B			00		
Antimony	ND	0.56	mg/kg	12/08/06	BSZ
Arsenic	ND	0.56	mg/kg	12/08/06	BSZ
Barium	14	0.56	mg/kg	12/08/06	BSZ
Beryllium	0.24	0.056	mg/kg	12/08/06	BSZ
Cadmium	0.15	0.11	mg/kg	12/08/06	BSZ
Chromium	6.0	0.56	mg/kg	12/08/06	BSZ
Copper	2.0	0.56	mg/kg	12/08/06	BSZ
Lead	3.4	0.23	mg/kg	12/11/06	BSZ
Nickel	2.0	0.56	mg/kg	12/08/06	BSZ
Selenium	ND	0.56	mg/kg	12/08/06	BSZ
Silver	ND	0.11	mg/kg	12/08/06	BSZ
Thallium	ND	0.28	mg/kg	12/08/06	BSZ
Zinc	4.0	0.56	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW	ND	0.023	mg/kg	12/05/06	AM

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 12/1/2006 Project: 20050458.B10/Nu-Style Phase II

Parameter		Result	DL	Units	Completed	By Dilution
(16) 841061130-16						
Date Collected: 11/30/2006	Matrix: Solid					
Cyanide, Total, by SW-846 9012	THE POIL	ND	0.53	mg/kg	12/06/06	DDD
Trace Metals by 6010B		1,2	0.00		12,00,00	222
Antimony		ND	0.53	mg/kg	12/08/06	BSZ
Arsenic		ND	0.53	mg/kg	12/08/06	BSZ
Barium		9.2	0.53	mg/kg	12/08/06	BSZ
Beryllium		0.081	0.053	mg/kg	12/08/06	BSZ
Cadmium		ND	0.11	mg/kg	12/08/06	BSZ
Chromium		1.4	0.53	mg/kg	12/08/06	BSZ
Copper		2.5	0.53	mg/kg	12/08/06	BSZ
Lead		1.5	0.21	mg/kg	12/11/06	BSZ
Nickel		6.5	0.53	mg/kg	12/08/06	BSZ
Selenium		ND	0.53	mg/kg	12/08/06	BSZ
Silver		ND	0.11	mg/kg	12/08/06	BSZ
Thallium		ND	0.27	mg/kg	12/08/06	BSZ
Zinc		4.2	0.53	mg/kg	12/08/06	BSZ
Mercury by SW-846 7471 in SW		ND	0.021	mg/kg	12/05/06	AM
(17) 841061130-17						
Date Collected: 11/30/2006	Matrix: Solid					
Cyanide, Total, by SW-846 9012		ND	0.55	mg/kg	12/06/06	DDD
Trace Metals by 6010B						
Antimony		ND	0.55	mg/kg	12/08/06	BSZ
Arsenic		ND	0.55	mg/kg	12/08/06	BSZ
Barium		17	0.55	mg/kg	12/08/06	BSZ
Beryllium		0.15	0.055	mg/kg	12/08/06	BSZ
Cadmium		ND	0.11	mg/kg	12/08/06	BSZ
Chromium		5.7	0.55	mg/kg	12/08/06	BSZ
Copper		25	0.55	mg/kg	12/08/06	BSZ
Lead		4.7	0.22	mg/kg	12/11/06	BSZ
Nickel		2.0	0.55	mg/kg	12/08/06	BSZ
Selenium		ND	0.55	mg/kg	12/08/06	BSZ
Silver		ND	0.11	mg/kg	12/08/06	BSZ
Thallium		ND	0.28	mg/kg	12/08/06	BSZ
Zinc		16	0.55	mg/kg	12/08/06	BSZ

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Laboratory: Premier Laboratory, LLC

PL Report No: E612052

Date Received: 12/1/2006

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II

 Parameter
 Result
 DL
 Units
 Completed
 By Dilution

 (17) 841061130-17 (continued)
 Date Collected: 11/30/2006

 Mercury by SW-846 7471 in SW
 ND
 0.022
 mg/kg
 12/05/06
 AM

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061130-01
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	5.8
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6200	ug/kg
C9-C12 Aliphatics**	50	ND	6200	ug/kg
C9-C10 Aromatics***	50	ND	6200	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	102	70%-130%
2,5-dibromotoluene #2	111	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	310	ug/kg
Ethylbenzene	ND	310	ug/kg
Methyl tert-butyl ether (MTBE)	ND	62	ug/kg
Naphthalene	ND	310	ug/kg
Toluene	ND	310	ug/kg
m,p-Xylenes	ND	310	ug/kg
o-Xylene	ND	310	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 Sample Description: 841061130-01

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:5.8Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28531.D;J28747.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	21
71-43-2	Benzene	ND	5.3
108-86-1	Bromobenzene	ND	5.3
74-97-5	Bromochloromethane	ND	5.3
75-27-4	Bromodichloromethane	ND	5.3
75-25-2	Bromoform	ND	5.3
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.3
135-98-8	sec-Butylbenzene	ND	5.3
98-06-6	tert-Butylbenzene	ND	5.3
75-15-0	Carbon disulfide	ND	5.3
56-23-5	Carbon tetrachloride	ND	5.3
108-90-7	Chlorobenzene	ND	5.3
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.3
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.3
106-43-4	4-Chlorotoluene	ND	5.3
108-20-3	Di-isopropyl ether (DIPE)	ND	53
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.3
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
74-95-3	Dibromomethane	ND	5.3
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.3
106-46-7	1,4-Dichlorobenzene	ND	5.3
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.3
107-06-2	1,2-Dichloroethane	ND	5.3
75-35-4	1,1-Dichloroethene	ND	5.3
156-59-2	cis-1,2-Dichloroethene	ND	5.3
156-60-5	trans-1,2-Dichloroethene	ND	5.3
78-87-5	1,2-Dichloropropane	ND	5.3
142-28-9	1,3-Dichloropropane	ND	5.3
590-20-7	2,2-Dichloropropane	ND	5.3
563-58-6	1,1-Dichloropropene	ND	5.3
10061-01-5	cis-1,3-Dichloropropene	ND	5.3
10061-02-6	trans-1,3-Dichloropropene	ND	5.3
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	21
/ -	.,	112	

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 (continued) Sample Description: 841061130-01

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 5.8

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50075 Lab Data File: J28531.D;J28747.D

Units: ug/kg

Ethyl tertiary-butyl ether 100-41-4 Ethylbenzene 87-68-3 Hexachlorobutadiene 591-78-6 2-Hexanone 1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone	er (EtBE)		ND	52
87-68-3 Hexachlorobutadiene 591-78-6 2-Hexanone 98-82-8 Isopropylbenzene 99-87-6 4-Isopropyltoluene 1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone				53
591-78-6 2-Hexanone 98-82-8 Isopropylbenzene 99-87-6 4-Isopropyltoluene 1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone			ND	5.3
98-82-8 Isopropylbenzene 99-87-6 4-Isopropyltoluene 1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone	Hexachlorobutadiene			5.3
99-87-6 4-Isopropyltoluene 1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone			ND	10
1634-04-4 Methyl tert-butyl ether 108-10-1 4-Methyl-2-pentanone			ND	5.3
108-10-1 4-Methyl-2-pentanone			ND	5.3
7 1	(MTBE)		ND	5.3
	(MIBK)		ND	10
75-09-2 Methylene chloride			ND	5.3
91-20-3 Naphthalene			ND	5.3
103-65-1 n-Propylbenzene			ND	5.3
100-42-5 Styrene			ND	5.3
994-05-8 Tertiary-amyl methyl e	ther (TAME)		ND	53
109-99-9 Tetrahydrofuran				5.3
96-18-4 1,2,3-Trichloropropane	-		ND	5.3
	1,1,1,2-Tetrachloroethane		ND	5.3
79-34-5 1,1,2,2-Tetrachloroetha	1,1,2,2-Tetrachloroethane		ND	5.0
127-18-4 Tetrachloroethene (PC)	Tetrachloroethene (PCE)		ND	5.3
108-88-3 Toluene	• • •		ND	5.3
87-61-6 1,2,3-Trichlorobenzene	1,2,3-Trichlorobenzene		ND	5.3
120-82-1 1,2,4-Trichlorobenzene	1,2,4-Trichlorobenzene		ND	5.3
71-55-6 1,1,1-Trichloroethane			ND	5.3
79-00-5 1,1,2-Trichloroethane			ND	5.3
79-01-6 Trichloroethene (TCE)			ND	5.3
75-69-4 Trichlorofluoromethan	ie		ND	10
95-63-6 1,2,4-Trimethylbenzene	e		ND	5.3
108-67-8 1,3,5-Trimethylbenzene	e		ND	5.3
75-01-4 Vinyl chloride			ND	10
95-47-6 o-Xylene			ND	5.3
m,p-Xylenes			ND	5.3
Surrogate	Recovery	Limits		
Bromofluorobenzene	86%	78%-111%		
1,2-Dichloroethane-d4	102%	91%-114%		
Toluene-d8	102%	86%-115%		

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061130-02
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	11.0
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6500	ug/kg
C9-C12 Aliphatics**	50	ND	6500	ug/kg
C9-C10 Aromatics***	50	ND	6500	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	98	70%-130%
2,5-dibromotoluene #2	108	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	330	ug/kg
Ethylbenzene	ND	330	ug/kg
Methyl tert-butyl ether (MTBE)	ND	65	ug/kg
Naphthalene	ND	330	ug/kg
Toluene	ND	330	ug/kg
m,p-Xylenes	ND	330	ug/kg
o-Xylene	ND	330	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 Sample Description: 841061130-02

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:11.0Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28532.D;J28748.D

Units: ug/kg

Actionary Acti	CAS No.	Parameter	Result	DL
108.86-1 Bromochloromethane ND 5.6 74-97-5 Bromochloromethane ND 5.6 74-97-5 Bromochloromethane ND 5.6 75-27-4 Bromofichloromethane ND 5.6 75-27-2 Bromofichloromethane ND 5.6 75-27-2 Bromoform ND 5.6 74-83-9 Bromomethane ND 11 74-83-3 2-Butanone (MEK) ND 11 74-83-3 2-Butanone (MEK) ND 11 74-97-3 74-	67-64-1	Acetone	ND	22
74-97-5 Bromochloromethane ND 5.6 75-27-4 Bromodichloromethane ND 5.6 75-25-2 Bromoform ND 5.6 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 78-93-3 2-Butanone (MEK) ND 15.6 104-51-8 n-Butylbenzene ND 5.6 135-98-8 sec-Butylbenzene ND 5.6 98-06-6 tert-Butylbenzene ND 5.6 98-06-6 tert-Butylbenzene ND 5.6 75-15-0 Carbon disulfide ND 5.6 56-23-5 Carbon tetrachloride ND 5.6 108-90-7 Chloroform ND 5.6 75-00-3 Chloroform ND 5.6 75-48-8 2-Chlorofolmen ND 5.6 108-90-8 2-Chlorotoluene ND 5.6 108-20-3 Di-sopropyl ether (DIPE) ND 5.6 108-	71-43-2	Benzene	ND	5.6
75-27-4 Bromofcom ND 5.6 75-25-2 Bromoform ND 5.6 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.6 135-98-8 sec-Butylbenzene ND 5.6 98-06-6 tert-Butylbenzene ND 5.6 98-06-6 tert-Butylbenzene ND 5.6 75-15-0 Carbon disulfide ND 5.6 66-23-5 Carbon detrachloride ND 5.6 108-90-7 Chlorobenzene ND 5.6 75-00-3 Chlorocethane ND 11 67-66-3 Chloroform ND 5.6 74-87-3 Chloromethane ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 108-20-3<	108-86-1	Bromobenzene	ND	5.6
75-25-2 Bromoferm	74-97-5	Bromochloromethane	ND	5.6
Page	75-27-4	Bromodichloromethane	ND	5.6
78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.6 135-98-8 sec-Butylbenzene ND 5.6 98-06-6 tert-Butylbenzene ND 5.6 75-15-0 Carbon disulfide ND 5.6 56-23-5 Carbon tetrachloride ND 5.6 108-90-7 Chlorobenzene ND 5.6 75-00-3 Chlorothane ND 11 67-66-3 Chloroform ND 5.6 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.6 104-34-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 106-34-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 12-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dichlorobenzene ND 5.6	75-25-2	Bromoform	ND	5.6
104-51-8	74-83-9	Bromomethane	ND	11
135-98-8 sec-Butylbenzene	78-93-3	2-Butanone (MEK)	ND	11
135-98-8 sec-Butylbenzene	104-51-8	n-Butylbenzene	ND	5.6
98-06-6 tert-Butylbenzene ND 5.6 75-15-0 Carbon disulfide ND 5.6 75-15-0 Carbon tetrachloride ND 5.6 108-90-7 Chlorobenzene ND 5.6 75-00-3 Chlorochane ND 11 67-66-3 Chloroform ND 5.6 74-87-3 Chlorochane ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 106-93-4 1,2-Dibromo-dhane (EDB) ND 5.6 106-93-4 1,2-Dibromochloromethane ND 5.0 95-50-1 1,2-Dichlorobenzene ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 107-06-2 1,2-Dichlorodifluoromethane ND 5.6 155-59-2 cis-1,2-Dichlorothene N	135-98-8	•	ND	5.6
56-23-5 Carbon tetrachloride ND 5.6 108-90-7 Chlorobenzene ND 5.6 75-00-3 Chlorotehane ND 11 67-66-3 Chloroform ND 5.6 74-87-3 Chlorotoluene ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 56 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 106-93-4 1,2-Dibromo-dhane (EDB) ND 5.6 106-93-4 1,2-Dibromochtane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 95-51-1 1,3-Dichlorobenzene ND 5.6 95-51-18 Dichlorodifluoromethane ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodhane ND	98-06-6	•	ND	5.6
108-90-7	75-15-0	Carbon disulfide	ND	5.6
75-00-3 Chloroethane ND 11 67-66-3 Chloroform ND 5.6 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.6 106-93-4 1,2-Dibromochlare (EDB) ND 3.4 74-95-3 Dibromomethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroptonethane ND <td>56-23-5</td> <td>Carbon tetrachloride</td> <td>ND</td> <td>5.6</td>	56-23-5	Carbon tetrachloride	ND	5.6
67-66-3 Chloroform ND 5.6 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 5.6 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 5.6 107-06-2 1,2-Dichlorothane ND 5.6 107-06-2 1,2-Dichlorothene ND 5.6 156-59-2 cis-1,2-Dichlorothene ND 5.6 156-60-5 trans-1,2-Dichloropropane	108-90-7	Chlorobenzene	ND	5.6
74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 56 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 1.6 107-06-2 1,2-Dichlorotethane ND 5.6 107-06-2 1,2-Dichlorotethane ND 5.6 156-59-2 cis-1,2-Dichlorotethene ND 5.6 156-60-5 trans-1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloroprop	75-00-3	Chloroethane	ND	11
95-49-8 2-Chlorotoluene ND 5.6 106-43-4 4-Chlorotoluene ND 5.6 108-20-3 Di-isopropyl ether (DIPE) ND 56 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloropropane ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 590-20-7 2,2-Dichloroprop	67-66-3	Chloroform	ND	5.6
106-43-4	74-87-3	Chloromethane	ND	11
108-20-3 Di-isopropyl ether (DIPE) ND 56 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-02-6 t	95-49-8	2-Chlorotoluene	ND	5.6
96-12-8 1,2-Dibrono-3-chloropropane (DBCP) ND 5.6 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.6 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 5.6 107-06-2 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 503-58-6 1,1-Dichloropropane ND 5.6 10061-02-6 trans-1	106-43-4	4-Chlorotoluene	ND	5.6
124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 503-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloro	108-20-3	Di-isopropyl ether (DIPE)	ND	56
106-93-4 1,2-Dibromoethane (EDB) ND 3.4 74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropane ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropr	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.6
74-95-3 Dibromomethane ND 5.6 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	124-48-1	Dibromochloromethane	ND	5.0
95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	106-93-4	1,2-Dibromoethane (EDB)	ND	3.4
541-73-1 1,3-Dichlorobenzene ND 5.6 106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	74-95-3	Dibromomethane	ND	5.6
106-46-7 1,4-Dichlorobenzene ND 5.6 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	95-50-1	1,2-Dichlorobenzene	ND	5.0
75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	541-73-1	1,3-Dichlorobenzene	ND	5.6
75-34-3 1,1-Dichloroethane ND 5.6 107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	106-46-7	1,4-Dichlorobenzene	ND	5.6
107-06-2 1,2-Dichloroethane ND 5.6 75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	75-71-8	Dichlorodifluoromethane	ND	11
75-35-4 1,1-Dichloroethene ND 5.6 156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	75-34-3	1,1-Dichloroethane	ND	5.6
156-59-2 cis-1,2-Dichloroethene ND 5.6 156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	107-06-2	1,2-Dichloroethane	ND	5.6
156-60-5 trans-1,2-Dichloroethene ND 5.6 78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	75-35-4	1,1-Dichloroethene	ND	5.6
78-87-5 1,2-Dichloropropane ND 5.6 142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	156-59-2	cis-1,2-Dichloroethene	ND	5.6
142-28-9 1,3-Dichloropropane ND 5.6 590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	156-60-5	trans-1,2-Dichloroethene	ND	5.6
590-20-7 2,2-Dichloropropane ND 5.6 563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	78-87-5	1,2-Dichloropropane	ND	5.6
563-58-6 1,1-Dichloropropene ND 5.6 10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	142-28-9	1,3-Dichloropropane	ND	5.6
10061-01-5 cis-1,3-Dichloropropene ND 5.6 10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	590-20-7	2,2-Dichloropropane	ND	5.6
10061-02-6 trans-1,3-Dichloropropene ND 5.6 60-29-7 Diethyl ether ND 11	563-58-6	1,1-Dichloropropene	ND	5.6
60-29-7 Diethyl ether ND 11	10061-01-5	cis-1,3-Dichloropropene	ND	5.6
·		trans-1,3-Dichloropropene	ND	5.6
123-91-1 1,4-Dioxane ND 22	60-29-7	Diethyl ether	ND	11
	123-91-1	1,4-Dioxane	ND	22

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Matrix: Solid

Percent Moisture: 11.0 Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 (continued) Sample Description: 841061130-02

Date Collected: 11/30/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/05/06 By: GP

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1
Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28532.D;J28748.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	56
100-41-4	Ethylbenzene	ND	5.6
87-68-3	Hexachlorobutadiene	ND	5.6
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.6
99-87-6	4-Isopropyltoluene	ND	5.6
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.6
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.6
91-20-3	Naphthalene	ND	5.6
103-65-1	n-Propylbenzene	ND	5.6
100-42-5	Styrene	ND	5.6
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	56
109-99-9	Tetrahydrofuran	ND	5.6
96-18-4	1,2,3-Trichloropropane	ND	5.6
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.6
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	ND	5.6
108-88-3	Toluene	ND	5.6
87-61-6	1,2,3-Trichlorobenzene	ND	5.6
120-82-1	1,2,4-Trichlorobenzene	ND	5.6
71-55-6	1,1,1-Trichloroethane	ND	5.6
79-00-5	1,1,2-Trichloroethane	ND	5.6
79-01-6	Trichloroethene (TCE)	ND	5.6
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.6
108-67-8	1,3,5-Trimethylbenzene	ND	5.6
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.6
	m,p-Xylenes	ND	5.6
Surrogate	Recovery Lir	nits	
Bromofluorobenzene	97% 78	%-111%	

 Surrogate
 Recovery
 Limits

 Bromofluorobenzene
 97%
 78%-111%

 1,2-Dichloroethane-d4
 102%
 91%-114%

 Toluene-d8
 94%
 86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061130-03
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	10.7
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6900	ug/kg
C9-C12 Aliphatics**	50	ND	6900	ug/kg
C9-C10 Aromatics***	50	ND	6900	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	96	70%-130%
2,5-dibromotoluene #2	105	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	350	ug/kg
Ethylbenzene	ND	350	ug/kg
Methyl tert-butyl ether (MTBE)	ND	69	ug/kg
Naphthalene	ND	350	ug/kg
Toluene	ND	350	ug/kg
m,p-Xylenes	ND	350	ug/kg
o-Xylene	ND	350	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 Sample Description: 841061130-03

Date Collected: 11/30/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Analyzed: 12/05/06 By: GP

Method: 8260B QC Batch#: 50075 Units: ug/kg Matrix: Solid Percent Moisture: 10.7 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume:

Lab Data File: J28533.D;J28749.D

Acetone	CAS No.	Parameter	Result	DL
108-86-1 Bromochloromethane ND 5.1 74-97-5 Bromochloromethane ND 5.1 75-27-4 Bromochloromethane ND 5.1 75-27-2 Bromoform ND 5.1 75-25-2 Bromoform ND 5.1 74-83-9 Bromomethane ND 10 74-83-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 156-23-5 Carbon disulfide ND 5.1 108-90-7 Chlorobenzene ND 5.1 108-90-7 Chlorobenzene ND 5.1 108-90-7 Chlorobenzene ND 5.1 108-90-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 104-95-3 Dibromomethane ND 5.1 105-90-1 1,2-Dichlorobenzene ND 5.1 106-93-4 1,2-Dibromomethane ND 5.1 106-93-4 1,2-Dichlorobenzene ND 5.1 106-93-4 1,2-Dichlorobenzene ND 5.1 106-93-4 1,2-Dichlorobenzene ND 5.1 106-95-2 1,2-Dichlorobenzene ND 5.1 106-96-2 1,2-Dichlorobenzene ND 5.1 107-06-2 1,2-Dichlorobenzene ND 5.1 156-60-5 trans-1,2-Dichloropenae ND 5.1 156-99-20-7 2,2-Dichloropenae ND 5.1 10061-02-6 trans-1,2-Dichloropene ND 5.1 10061-02-6	67-64-1	Acetone	ND	20
74-97-5 Bromochloromethane ND 5.1 75-27-4 Bromodichloromethane ND 5.1 75-25-2 Bromoform ND 5.1 74-83-9 Bromomethane ND 10 78-93-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 98-06-6 tert-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 56-23-5 Carbon tetrachloride ND 5.1 108-90-7 Chlorobenzene ND 5.1 75-00-3 Chlorocethane ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chlorocethane ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 12-48-1 Dibromochloromethane ND 5.0	71-43-2	Benzene	ND	5.1
75-27-4 Bromoform ND 5.1 75-25-2 Bromoform ND 5.1 74-83-9 Bromomethane ND 10 78-93-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 185-98-8 sec-Butylbenzene ND 5.1 98-06-6 tert-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 108-90-7 Chloroben tetrachloride ND 5.1 108-90-7 Chlorobenzene ND 5.1 75-00-3 Chlorobenzene ND 5.1 75-00-3 Chlorobentene ND 5.1 74-87-3 Chlorotoluene ND 5.1 98-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 105-20-3 Di-isopropyl ether (DIPE) ND 5.1 96-12-8	108-86-1	Bromobenzene	ND	5.1
75-25-2 Bromoform ND 5.1 74-83-9 Bromomethane ND 10 78-93-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 75-15-0 Carbon tetrachloride ND 5.1 108-90-7 Chlorotenzene ND 5.1 75-00-3 Chlorotene ND 5.1 75-00-3 Chlorotene ND 5.1 74-87-3 Chlorotene ND 5.1 74-87-3 Chloroteluene ND 5.1 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 106-3-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 106-93-4 1,2-Dibromo-dhoromethane ND 5.0 142-8	74-97-5	Bromochloromethane	ND	5.1
74-83-9 Bromomethane ND 10 78-93-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 98-06-6 tert-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 56-23-5 Carbon tetrachloride ND 5.1 108-90-7 Chlorofene ND 5.1 75-00-3 Chlorofene ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chlorofoluene ND 5.1 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-joromo-s-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 109-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1	75-27-4	Bromodichloromethane	ND	5.1
78-93-3 2-Butanone (MEK) ND 10 104-51-8 n-Butylbenzene ND 5.1 135-98-8 sec-Butylbenzene ND 5.1 98-06-6 tert-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 56-23-5 Carbon tetrachloride ND 5.1 108-90-7 Chlorobenzene ND 5.1 75-00-3 Chlorothane ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chloromethane ND 5.1 95-49-8 2-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 106-93-4 1,2-Dirbomoethane (EDB) ND 5.1 106-93-4 1,2-Dirbomoethane (EDB) ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.1 </td <td>75-25-2</td> <td>Bromoform</td> <td>ND</td> <td>5.1</td>	75-25-2	Bromoform	ND	5.1
104-51-8	74-83-9	Bromomethane	ND	10
135-98-8 sec-Butylbenzene	78-93-3	2-Butanone (MEK)	ND	10
135-98-8 sec-Butylbenzene	104-51-8	n-Butylbenzene	ND	5.1
98-06-6 tert-Butylbenzene ND 5.1 75-15-0 Carbon disulfide ND 5.1 75-15-0 Carbon disulfide ND 5.1 108-90-7 Chlorobenzene ND 5.1 75-00-3 Chlorotehane ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chlorotehane ND 5.1 95-49-8 2-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE)	135-98-8		ND	5.1
75-15-0 Carbon disulfide ND 5.1 56-23-5 Carbon tetrachloride ND 5.1 108-90-7 Chlorobenzene ND 5.1 75-00-3 Chlorochtane ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chlorotoluene ND 5.1 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 104-48-1 Dibromo-s-chloropropane (DBCP) ND 5.1 124-48-1 Dibromo-chloromethane ND 5.0 106-93-4 1,2-Dibromo-s-chloropropane (DBCP) ND 5.1 14-95-3 Dibromo-chloromethane ND 5.1 74-95-3 Dibromo-chlorobenzene ND 5.1 106-46-7 1,2-Dichlorobenzene ND 5.1 5-17-7-18 Dichlorodifluoromethane	98-06-6	· · · · · · · · · · · · · · · · · · ·	ND	5.1
108-90-7	75-15-0	· · · · · · · · · · · · · · · · · · ·	ND	5.1
75-00-3 Chloroethane ND 10 67-66-3 Chloroform ND 5.1 74-87-3 Chloromethane ND 10 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 51 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 106-93-4 1,2-Dibromoethane (EDB) ND 5.1 195-50-1 1,2-Dichlorobenzene ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.1 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethene	56-23-5	Carbon tetrachloride	ND	5.1
67-66-3 Chloroform ND 5.1 74-87-3 Chloromethane ND 10 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 5.1 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochloromethane ND 5.1 74-95-3 Dibromochloromethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethane	108-90-7	Chlorobenzene	ND	5.1
74-87-3 Chloromethane ND 10 95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 51 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.1 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 107-06-2 1,2-Dichlorobenzene ND 5.1 107-06-2 1,2-Dichlorothane ND 5.1 156-59-2 cis-1,2-Dichlorothene ND 5.1 156-60-5 trans-1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane<	75-00-3	Chloroethane	ND	10
95-49-8 2-Chlorotoluene ND 5.1 106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 51 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloropropane ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloro	67-66-3	Chloroform	ND	5.1
106-43-4 4-Chlorotoluene ND 5.1 108-20-3 Di-isopropyl ether (DIPE) ND 51 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 75-34-3 1,1-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 563-58-6 1,1-Dichlo	74-87-3	Chloromethane	ND	10
108-20-3 Di-isopropyl ether (DIPE) ND 51 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Di	95-49-8	2-Chlorotoluene	ND	5.1
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 107-06-2 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6	106-43-4	4-Chlorotoluene	ND	5.1
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.1 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 5.1 107-06-2 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6	108-20-3	Di-isopropyl ether (DIPE)	ND	51
124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroptopethene ND 5.1 18-78-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropane ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichl	96-12-8	* **	ND	5.1
106-93-4 1,2-Dibromoethane (EDB) ND 3.1 74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 503-58-6 1,1-Dichloropropane ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	124-48-1	* * * * * * * * * * * * * * * * * * *	ND	5.0
74-95-3 Dibromomethane ND 5.1 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 5.1	106-93-4		ND	3.1
541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	74-95-3		ND	5.1
541-73-1 1,3-Dichlorobenzene ND 5.1 106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	95-50-1	1,2-Dichlorobenzene	ND	5.0
106-46-7 1,4-Dichlorobenzene ND 5.1 75-71-8 Dichlorodifluoromethane ND 10 75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	541-73-1		ND	5.1
75-34-3 1,1-Dichloroethane ND 5.1 107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	106-46-7		ND	5.1
107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	75-71-8	Dichlorodifluoromethane	ND	10
107-06-2 1,2-Dichloroethane ND 5.1 75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	75-34-3	1,1-Dichloroethane	ND	5.1
75-35-4 1,1-Dichloroethene ND 5.1 156-59-2 cis-1,2-Dichloroethene ND 5.1 156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	107-06-2		ND	5.1
156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	75-35-4		ND	5.1
156-60-5 trans-1,2-Dichloroethene ND 5.1 78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	156-59-2	cis-1,2-Dichloroethene	ND	5.1
78-87-5 1,2-Dichloropropane ND 5.1 142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	156-60-5	trans-1,2-Dichloroethene	ND	5.1
142-28-9 1,3-Dichloropropane ND 5.1 590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	78-87-5	1,2-Dichloropropane	ND	5.1
590-20-7 2,2-Dichloropropane ND 5.1 563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	142-28-9		ND	5.1
563-58-6 1,1-Dichloropropene ND 5.1 10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10	590-20-7		ND	5.1
10061-01-5 cis-1,3-Dichloropropene ND 5.1 10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10				
10061-02-6 trans-1,3-Dichloropropene ND 5.1 60-29-7 Diethyl ether ND 10				
60-29-7 Diethyl ether ND 10				
·				
		1,4-Dioxane		

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Customer: Fuss & O'Neill Laboratory: Premier Laboratory, LLC Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 (continued) Sample Description: 841061130-03

Date Collected: 11/30/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 10.7 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/05/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume: Lab Data File: J28533.D;J28749.D QC Batch#: 50075

Units: ug/kg

CAS No.	Parameter	Result	DL		
	Ethyl tertiary-butyl ether (EtBE)	ND	51		
100-41-4	Ethylbenzene	ND	5.1		
87-68-3	Hexachlorobutadiene	ND	5.1		
591-78-6	2-Hexanone	ND	10		
98-82-8	Isopropylbenzene	ND	5.1		
99-87-6	4-Isopropyltoluene	ND	5.1		
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.1		
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10		
75-09-2	Methylene chloride	ND	5.1		
91-20-3	Naphthalene	ND	5.1		
103-65-1	n-Propylbenzene	ND	5.1		
100-42-5	Styrene	ND	5.1		
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	51		
109-99-9	Tetrahydrofuran	ND	5.1		
96-18-4	1,2,3-Trichloropropane	ND	5.1		
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.1		
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0		
127-18-4	Tetrachloroethene (PCE)	ND	5.1		
108-88-3	Toluene	17	5.1		
87-61-6	1,2,3-Trichlorobenzene	ND	5.1		
120-82-1	1,2,4-Trichlorobenzene				
71-55-6	1,1,1-Trichloroethane				
79-00-5	1,1,2-Trichloroethane	ND	5.1		
79-01-6	Trichloroethene (TCE)	ND	5.1		
75-69-4	Trichlorofluoromethane	ND	10		
95-63-6	1,2,4-Trimethylbenzene	ND	5.1		
108-67-8	1,3,5-Trimethylbenzene	ND	5.1		
75-01-4	Vinyl chloride	ND	10		
95-47-6	o-Xylene	ND	5.1		
	m,p-Xylenes	7.0	5.1		
Surrogate	Recovery	Limits			
Bromofluorobenzene	89%	78%-111%			
1,2-Dichloroethane-d4	103%	91%-114%			
Toluono do	1010/	960/ 1150/			

Toluene-d8 101% 86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061130-04
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	15.8
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	8800	ug/kg
C9-C12 Aliphatics**	50	ND	8800	ug/kg
C9-C10 Aromatics***	50	ND	8800	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	105	70%-130%
2,5-dibromotoluene #2	112	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	440	ug/kg
Ethylbenzene	ND	440	ug/kg
Methyl tert-butyl ether (MTBE)	ND	88	ug/kg
Naphthalene	ND	440	ug/kg
Toluene	ND	440	ug/kg
m,p-Xylenes	ND	440	ug/kg
o-Xylene	ND	440	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 Sample Description: 841061130-04

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:15.8Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/04/06By:Dilution Factor:1

Method: 8260B Soil Extract Volume:
QC Batch#: 50061 Lab Data File: J28510.D;J28750.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	23
71-43-2	Benzene	ND	5.8
108-86-1	Bromobenzene	ND	5.8
74-97-5	Bromochloromethane	ND	5.8
75-27-4	Bromodichloromethane	ND	5.8
75-25-2	Bromoform	ND	5.8
74-83-9	Bromomethane	ND	12
78-93-3	2-Butanone (MEK)	ND	12
104-51-8	n-Butylbenzene	ND	5.8
135-98-8	sec-Butylbenzene	ND	5.8
98-06-6	tert-Butylbenzene	ND	5.8
75-15-0	Carbon disulfide	ND	5.8
56-23-5	Carbon tetrachloride	ND	5.8
108-90-7	Chlorobenzene	ND	5.8
75-00-3	Chloroethane	ND	12
67-66-3	Chloroform	ND	5.8
74-87-3	Chloromethane	ND	12
95-49-8	2-Chlorotoluene	ND	5.8
106-43-4	4-Chlorotoluene	ND	5.8
108-20-3	Di-isopropyl ether (DIPE)	ND	58
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.8
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.5
74-95-3	Dibromomethane	ND	5.8
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.8
106-46-7	1,4-Dichlorobenzene	ND	5.8
75-71-8	Dichlorodifluoromethane	ND	12
75-34-3	1,1-Dichloroethane	ND	5.8
107-06-2	1,2-Dichloroethane	ND	5.8
75-35-4	1,1-Dichloroethene	ND	5.8
156-59-2	cis-1,2-Dichloroethene	ND	5.8
156-60-5	trans-1,2-Dichloroethene	ND	5.8
78-87-5	1,2-Dichloropropane	ND	5.8
142-28-9	1,3-Dichloropropane	ND	5.8
590-20-7	2,2-Dichloropropane	ND	5.8
563-58-6	1,1-Dichloropropene	ND	5.8
10061-01-5	cis-1,3-Dichloropropene	ND	5.8
10061-02-6	trans-1,3-Dichloropropene	ND	5.8
60-29-7	Diethyl ether	ND	12
123-91-1	1,4-Dioxane	ND	23

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 (continued) Sample Description: 841061130-04

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 15.8
Date Extracted: By: Sample Weight/Volume:
Date Analyzed: 12/04/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:
QC Batch#: 50061 Lab Data File: J28510.D;J28750.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	58
100-41-4	Ethylbenzene	ND	5.8
87-68-3	Hexachlorobutadiene	ND	5.8
591-78-6	2-Hexanone	ND	12
98-82-8	Isopropylbenzene	ND	5.8
99-87-6	4-Isopropyltoluene	ND	5.8
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.8
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	12
75-09-2	Methylene chloride	ND	5.8
91-20-3	Naphthalene	10	5.8
103-65-1	n-Propylbenzene	ND	5.8
100-42-5	Styrene	ND	5.8
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	58
109-99-9	Tetrahydrofuran	ND	5.8
96-18-4	1,2,3-Trichloropropane	ND	5.8
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.8
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	ND	5.8
108-88-3	Toluene	ND	5.8
87-61-6	1,2,3-Trichlorobenzene	ND	5.8
120-82-1	1,2,4-Trichlorobenzene	ND	5.8
71-55-6	1,1,1-Trichloroethane	ND	5.8
79-00-5	1,1,2-Trichloroethane	ND	5.8
79-01-6	Trichloroethene (TCE)	ND	5.8
75-69-4	Trichlorofluoromethane	ND	12
95-63-6	1,2,4-Trimethylbenzene	ND	5.8
108-67-8	1,3,5-Trimethylbenzene	ND	5.8
75-01-4	Vinyl chloride	ND	12
95-47-6	o-Xylene	ND	5.8
	m,p-Xylenes	ND	5.8
Surrogate	Recovery	Limits	
Bromofluorobenzene	96%	78%-111%	
1.2 Dichloroethane d/	103%	01% 11/1%	

 Bromofluorobenzene
 96%
 78%-111%

 1,2-Dichloroethane-d4
 103%
 91%-114%

 Toluene-d8
 97%
 86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061130-05
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	14.7
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7900	ug/kg
C9-C12 Aliphatics**	50	ND	7900	ug/kg
C9-C10 Aromatics***	50	ND	7900	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	107	70%-130%
2,5-dibromotoluene #2	112	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	390	ug/kg
Ethylbenzene	ND	390	ug/kg
Methyl tert-butyl ether (MTBE)	ND	79	ug/kg
Naphthalene	ND	390	ug/kg
Toluene	ND	390	ug/kg
m,p-Xylenes	ND	390	ug/kg
o-Xylene	ND	390	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841061130-05

Date Collected: 11/30/2006Matrix: SolidDate Received: 12/1/2006Percent Moisture: 14.7Date Extracted: By: Sample Weight/Volume: Date Analyzed: 12/04/06By: GPDilution Factor: 1

Method: 8260B Soil Extract Volume:
QC Batch#: 50061 Lab Data File: J28511.D;M32618.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	21
71-43-2	Benzene	ND	5.3
108-86-1	Bromobenzene	ND	5.3
74-97-5	Bromochloromethane	ND	5.3
75-27-4	Bromodichloromethane	ND	5.3
75-25-2	Bromoform	ND	5.3
74-83-9	Bromomethane	ND	11
78-93-3	2-Butanone (MEK)	ND	11
104-51-8	n-Butylbenzene	ND	5.3
135-98-8	sec-Butylbenzene	ND	5.3
98-06-6	tert-Butylbenzene	ND	5.3
75-15-0	Carbon disulfide	ND	5.3
56-23-5	Carbon tetrachloride	ND	5.3
108-90-7	Chlorobenzene	ND	5.3
75-00-3	Chloroethane	ND	11
67-66-3	Chloroform	ND	5.3
74-87-3	Chloromethane	ND	11
95-49-8	2-Chlorotoluene	ND	5.3
106-43-4	4-Chlorotoluene	ND	5.3
108-20-3	Di-isopropyl ether (DIPE)	ND	53
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.3
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
74-95-3	Dibromomethane	ND	5.3
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.3
106-46-7	1,4-Dichlorobenzene	ND	5.3
75-71-8	Dichlorodifluoromethane	ND	11
75-34-3	1,1-Dichloroethane	ND	5.3
107-06-2	1,2-Dichloroethane	ND	5.3
75-35-4	1,1-Dichloroethene	ND	5.3
156-59-2	cis-1,2-Dichloroethene	ND	5.3
156-60-5	trans-1,2-Dichloroethene	ND	5.3
78-87-5	1,2-Dichloropropane	ND	5.3
142-28-9	1,3-Dichloropropane	ND	5.3
590-20-7	2,2-Dichloropropane	ND	5.3
563-58-6	1,1-Dichloropropene	ND	5.3
10061-01-5	cis-1,3-Dichloropropene	ND	5.3
10061-02-6	trans-1,3-Dichloropropene	ND	5.3
60-29-7	Diethyl ether	ND	11
123-91-1	1,4-Dioxane	ND	21
	,		

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Matrix: Solid

Percent Moisture: 14.7 Sample Weight/Volume:

Dilution Factor: 1

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 (continued) Sample Description: 841061130-05

Date Collected: 11/30/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/04/06 By: GP

Method: 8260B Soil Extract Volume:
QC Batch#: 50061 Lab Data File: J28511.D;M32618.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	53
100-41-4	Ethylbenzene	ND	5.3
87-68-3	Hexachlorobutadiene	ND	5.3
591-78-6	2-Hexanone	2-Hexanone ND	
98-82-8	Isopropylbenzene	ND	5.3
99-87-6	4-Isopropyltoluene	ND	5.3
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.3
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.3
91-20-3	Naphthalene	ND	5.3
103-65-1	n-Propylbenzene	ND	5.3
100-42-5	Styrene	ND	5.3
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	53
109-99-9	Tetrahydrofuran	ND	5.3
96-18-4	1,2,3-Trichloropropane	ND	5.3
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.3
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	28	5.3
108-88-3	Toluene	ND	5.3
87-61-6	1,2,3-Trichlorobenzene	ND	5.3
120-82-1	1,2,4-Trichlorobenzene	ND	5.3
71-55-6	1,1,1-Trichloroethane ND		5.3
79-00-5	1,1,2-Trichloroethane	ND	5.3
79-01-6	Trichloroethene (TCE)	12	5.3
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.3
108-67-8	1,3,5-Trimethylbenzene	ND	5.3
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.3
	m,p-Xylenes	ND	5.3
Surrogate	Recovery Lin	mits	
Bromofluorobenzene	740/ 780/ 1110/		

 Surrogate
 Recovery
 Limits

 Bromofluorobenzene
 74%
 78%-111%

 1,2-Dichloroethane-d4
 101%
 91%-114%

 Toluene-d8
 115%
 86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061130-06
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	12.3
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7000	ug/kg
C9-C12 Aliphatics**	50	ND	7000	ug/kg
C9-C10 Aromatics***	50	ND	7000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	102	70%-130%
2,5-dibromotoluene #2	110	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	350	ug/kg
Ethylbenzene	ND	350	ug/kg
Methyl tert-butyl ether (MTBE)	ND	70	ug/kg
Naphthalene	ND	350	ug/kg
Toluene	ND	350	ug/kg
m,p-Xylenes	ND	350	ug/kg
o-Xylene	ND	350	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Matrix: Solid

Percent Moisture: 12.3

Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 6 Sample Description: 841061130-06

Date Collected: 11/30/2006

Date Received: 12/1/2006

Date Extracted: By:

Date Applying 12/05/06 By: GP

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1
Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28535.D;M32619.D

Units: ug/kg

Formation	CAS No.	Parameter	Result	DL
108-86-1 Bromobenzene	67-64-1	Acetone	ND	22
Page	71-43-2	Benzene	ND	5.5
75-27-4 Bromoform ND 5.5 75-25-2 Bromoform ND 5.5 74-83-9 Bromomethane ND 11 78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.5 135-98-8 sec-Butylbenzene ND 5.5 98-06-6 tert-Butylbenzene ND 5.5 75-15-0 Carbon disulfide ND 5.5 56-23-5 Carbon disulfide ND 5.5 108-90-7 Chlorobenzene ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chloroform ND 5.5 74-87-3 Chlorotoluene ND 5.5 106-43-4 4-Chiorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 106-48-1 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 <t< td=""><td>108-86-1</td><td>Bromobenzene</td><td>ND</td><td>5.5</td></t<>	108-86-1	Bromobenzene	ND	5.5
TS-25-2	74-97-5	Bromochloromethane	ND	5.5
A+83-9	75-27-4	Bromodichloromethane	ND	5.5
78-93-3 2-Butanone (MEK) ND 11 104-51-8 n-Butylbenzene ND 5.5 135-98-8 sec-Butylbenzene ND 5.5 98-06-6 tert-Butylbenzene ND 5.5 75-15-0 Carbon disulfide ND 5.5 56-23-5 Carbon tetrachloride ND 5.5 108-90-7 Chlorobenzene ND 5.5 75-00-3 Chlorotethane ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chlorotoluene ND 11 95-49-8 2-Chlorotoluene ND 5.5 104-34-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 106-93-4 1,2-Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochloromethane ND 5.5 94-1-73-1 1,3-Dichlorobenzene ND 5.	75-25-2	Bromoform	ND	5.5
104-51-8	74-83-9	Bromomethane	ND	11
135-98-8 sec-Dutylbenzene	78-93-3	2-Butanone (MEK)	ND	11
98-06-6 tert-Butylbenzene ND 5.5 75-15-0 Carbon disulfide ND 5.5 55-23-5 Carbon tetrachloride ND 5.5 108-90-7 Chlorobenzene ND 5.5 75-00-3 Chlorotethane ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chlorotethane ND 11 95-49-8 2-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 106-93-4 1,2-Dibromoedhane (EDB) ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromoethane (EDB) ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 96-12-8 1,2-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.0 15-5-91-8 Dichlorodifluoromethane ND <td>104-51-8</td> <td>n-Butylbenzene</td> <td>ND</td> <td>5.5</td>	104-51-8	n-Butylbenzene	ND	5.5
75-15-0 Carbon disulfide ND 5.5 56-23-5 Carbon tetrachloride ND 5.5 108-90-7 Chlorobenzene ND 5.5 75-00-3 Chlorobenzene ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chlorotoluene ND 11 95-49-8 2-Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 108-22-8 1,2-Di-isopropyl ether (DIPE) ND 5.5 106-29-3 1,2-Di-isopropyl ether	135-98-8	sec-Butylbenzene	ND	5.5
56-23-5 Carbon tetrachloride ND 5.5 108-90-7 Chlorobenzene ND 5.5 75-00-3 Chlorotehane ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chlorotehane ND 5.5 74-87-3 Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 104-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromochlane (EDB) ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 75-71-8 Dichlorodhane ND	98-06-6	tert-Butylbenzene	ND	5.5
108-90-7	75-15-0	Carbon disulfide	ND	5.5
75-00-3 Chloroethane ND 11 67-66-3 Chloroform ND 5.5 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 55 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 106-93-4 1,2-Dibromo-1-chloroptopane (DBCP) ND 5.0 106-93-4 1,2-Dibromo-thane (EDB) ND 5.0 95-50-1 1,2-Dichlorobenzene ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 107-06-2 1,2-Dichlorobenzene ND 5.5 107-06-2 1	56-23-5	Carbon tetrachloride	ND	5.5
67-66-3 Chloroform ND 5.5 74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloropropane ND 5.5 156-60-5 trans-1,2-Dichloropropane </td <td>108-90-7</td> <td>Chlorobenzene</td> <td>ND</td> <td>5.5</td>	108-90-7	Chlorobenzene	ND	5.5
74-87-3 Chloromethane ND 11 95-49-8 2-Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 55 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 55 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.5 541-73-1 1,3-Dichlorobenzene ND 5.5 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene	75-00-3	Chloroethane	ND	11
95-49-8 2-Chlorotoluene ND 5.5 106-43-4 4-Chlorotoluene ND 5.5 108-20-3 Di-isopropyl ether (DIPE) ND 5.5 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloropropane ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 590-20-7 2,2-Dichlor	67-66-3	Chloroform	ND	5.5
106-43-4	74-87-3	Chloromethane	ND	11
108-20-3 Di-isopropyl ether (DIPE) ND 55 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 c	95-49-8	2-Chlorotoluene	ND	5.5
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 5.5 124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloropropane ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropane ND 5.5 10061-01-5 cis-1,3-Di	106-43-4	4-Chlorotoluene	ND	5.5
124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 500-20-7 2,2-Dichloropropane ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloro	108-20-3	Di-isopropyl ether (DIPE)	ND	55
124-48-1 Dibromochloromethane ND 5.0 106-93-4 1,2-Dibromoethane (EDB) ND 3.3 74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 500-20-7 2,2-Dichloropropane ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloro	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.5
74-95-3 Dibromomethane ND 5.5 95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	124-48-1		ND	5.0
95-50-1 1,2-Dichlorobenzene ND 5.0 541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	106-93-4	1,2-Dibromoethane (EDB)	ND	3.3
541-73-1 1,3-Dichlorobenzene ND 5.5 106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	74-95-3	Dibromomethane	ND	5.5
106-46-7 1,4-Dichlorobenzene ND 5.5 75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	95-50-1	1,2-Dichlorobenzene	ND	5.0
75-71-8 Dichlorodifluoromethane ND 11 75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	541-73-1	1,3-Dichlorobenzene	ND	5.5
75-34-3 1,1-Dichloroethane ND 5.5 107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	106-46-7	1,4-Dichlorobenzene	ND	5.5
107-06-2 1,2-Dichloroethane ND 5.5 75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	75-71-8	Dichlorodifluoromethane	ND	11
75-35-4 1,1-Dichloroethene ND 5.5 156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	75-34-3	1,1-Dichloroethane	ND	5.5
156-59-2 cis-1,2-Dichloroethene ND 5.5 156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	107-06-2	1,2-Dichloroethane	ND	5.5
156-60-5 trans-1,2-Dichloroethene ND 5.5 78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	75-35-4	1,1-Dichloroethene	ND	5.5
78-87-5 1,2-Dichloropropane ND 5.5 142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	156-59-2	cis-1,2-Dichloroethene	ND	5.5
142-28-9 1,3-Dichloropropane ND 5.5 590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	156-60-5	trans-1,2-Dichloroethene	ND	5.5
590-20-7 2,2-Dichloropropane ND 5.5 563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	78-87-5	1,2-Dichloropropane	ND	5.5
563-58-6 1,1-Dichloropropene ND 5.5 10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	142-28-9	1,3-Dichloropropane	ND	5.5
10061-01-5 cis-1,3-Dichloropropene ND 5.5 10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	590-20-7	2,2-Dichloropropane	ND	5.5
10061-02-6 trans-1,3-Dichloropropene ND 5.5 60-29-7 Diethyl ether ND 11	563-58-6	1,1-Dichloropropene	ND	5.5
60-29-7 Diethyl ether ND 11	10061-01-5	cis-1,3-Dichloropropene	ND	5.5
· ·		trans-1,3-Dichloropropene	ND	5.5
123-91-1 1,4-Dioxane ND 22		Diethyl ether	ND	11
	123-91-1	1,4-Dioxane	ND	22

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Matrix: Solid

Percent Moisture: 12.3 Sample Weight/Volume:

Dilution Factor: 1

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 6 (continued) Sample Description: 841061130-06

Date Collected: 11/30/2006 Date Received: 12/1/2006 Date Extracted: By: Date Analyzed: 12/05/06 By: GP

Method: 8260B Soil Extract Volume: QC Batch#: 50075 Lab Data File: J28535.D;M32619.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	55
100-41-4	Ethylbenzene	ND	5.5
87-68-3	Hexachlorobutadiene	ND	5.5
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.5
99-87-6	4-Isopropyltoluene	ND	5.5
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.5
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.5
91-20-3	Naphthalene	ND	5.5
103-65-1	n-Propylbenzene	ND	5.5
100-42-5	Styrene	ND	5.5
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	55
109-99-9	Tetrahydrofuran	ND	5.5
96-18-4	1,2,3-Trichloropropane	ND	5.5
630-20-6	1,1,2-Tetrachloroethane	ND	5.5
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	45	5.5
108-88-3	Toluene	ND	5.5
87-61-6	1,2,3-Trichlorobenzene	ND	5.5
120-82-1	1,2,4-Trichlorobenzene	ND	5.5
71-55-6	1,1,1-Trichloroethane	ND	5.5
79-00-5	1,1,2-Trichloroethane	ND	5.5
79-01-6	Trichloroethene (TCE)	21	5.5
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.5
108-67-8	1,3,5-Trimethylbenzene	ND	5.5
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.5
	m,p-Xylenes	ND	5.5
Surrogate	Recovery	Limits	
Bromofluorobenzene	84%	78%-111%	
1.2 Dichloroethane d/	108%	01% 11/1%	

Surrogate	Recovery	Limits
Bromofluorobenzene	84%	78%-111%
1,2-Dichloroethane-d4	108%	91%-114%
Toluene-d8	104%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	7	Sample Description:	841061130-07
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	12.3
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7000	ug/kg
C9-C12 Aliphatics**	50	ND	7000	ug/kg
C9-C10 Aromatics***	50	ND	7000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	101	70%-130%
2,5-dibromotoluene #2	107	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	350	ug/kg
Ethylbenzene	ND	350	ug/kg
Methyl tert-butyl ether (MTBE)	ND	70	ug/kg
Naphthalene	ND	350	ug/kg
Toluene	ND	350	ug/kg
m,p-Xylenes	ND	350	ug/kg
o-Xylene	ND	350	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 Sample Description: 841061130-07

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 12.3
Date Extracted: By: Sample Weight/Volume

Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By: GPDilution Factor:100Method:8260BSoil Extract Volume:

QC Batch#: 51012 Lab Data File: M32519.D;M32620.D

Units: ug/kg

67-64-1 Acetone ND 2300 71-43-2 Benzene ND 570 108-86-1 Bromobenzene ND 570 74-97-5 Bromochloromethane ND 570 75-27-4 Bromodichloromethane ND 570 75-25-2 Bromoform ND 570 74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 75-15-0 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chlorotehane ND 570 74-87-3 Chlorotehane ND 570 74-88-3 2-Chlorotoluene ND 570 108-20-3 Di-
108-86-1 Bromobenzene ND 570 74-97-5 Bromochloromethane ND 570 75-27-4 Bromodichloromethane ND 570 75-25-2 Bromoform ND 570 74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chlorobenzene ND 570 74-87-3 Chloroform ND 570 74-87-3 Chloromethane ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
74-97-5 Bromochloromethane ND 570 75-27-4 Bromodichloromethane ND 570 75-25-2 Bromoform ND 570 74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 570 74-87-3 Chloroform ND 570 74-88-3 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
75-27-4 Bromodichloromethane ND 570 75-25-2 Bromoform ND 570 74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
75-25-2 Bromoform ND 570 74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
74-83-9 Bromomethane ND 1100 78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
78-93-3 2-Butanone (MEK) ND 1100 104-51-8 n-Butylbenzene ND 570 135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
135-98-8 sec-Butylbenzene ND 570 98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
98-06-6 tert-Butylbenzene ND 570 75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
75-15-0 Carbon disulfide ND 570 56-23-5 Carbon tetrachloride ND 570 108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
108-90-7 Chlorobenzene ND 570 75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
75-00-3 Chloroethane ND 1100 67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
67-66-3 Chloroform ND 570 74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
74-87-3 Chloromethane ND 1100 95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
95-49-8 2-Chlorotoluene ND 570 106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
106-43-4 4-Chlorotoluene ND 570 108-20-3 Di-isopropyl ether (DIPE) ND 57
Di-isopropyl ether (DIPE) ND 57
96.12.8 1.2-Dibromo_3_chloropropaga (DRCP) ND 570
70-12-0 1,2-DIOIOINO-3-CHIOIOPIOPAIIC (DDCF) ND 3/0
Dibromochloromethane ND 570
106-93-4 1,2-Dibromoethane (EDB) ND 340
74-95-3 Dibromomethane ND 570
95-50-1 1,2-Dichlorobenzene ND 570
541-73-1 1,3-Dichlorobenzene ND 570
106-46-7 1,4-Dichlorobenzene ND 570
75-71-8 Dichlorodifluoromethane ND 1100
75-34-3 1,1-Dichloroethane ND 570
107-06-2 1,2-Dichloroethane ND 570
75-35-4 1,1-Dichloroethene ND 570
156-59-2 cis-1,2-Dichloroethene ND 570
156-60-5 trans-1,2-Dichloroethene ND 570
78-87-5 1,2-Dichloropropane ND 570
142-28-9 1,3-Dichloropropane ND 570
590-20-7 2,2-Dichloropropane ND 570
563-58-6 1,1-Dichloropropene ND 570
10061-01-5 cis-1,3-Dichloropropene ND 570
10061-02-6 trans-1,3-Dichloropropene ND 570
60-29-7 Diethyl ether ND 1100
123-91-1 1,4-Dioxane ND 2300

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Matrix: Solid

Percent Moisture: 12.3 Sample Weight/Volume:

Dilution Factor: 100

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 7 (continued) Sample Description: 841061130-07

Date Collected: 11/30/2006 Date Received: 12/1/2006 Date Extracted: By: Date Analyzed: 12/06/06 By: GP

Method: 8260B Soil Extract Volume: QC Batch#: 51012 Lab Data File: M32519.D;M32620.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	57
100-41-4	Ethylbenzene	ND	570
87-68-3	Hexachlorobutadiene	ND	570
591-78-6	2-Hexanone	ND	1100
98-82-8	Isopropylbenzene	ND	570
99-87-6	4-Isopropyltoluene	ND	570
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	570
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	1100
75-09-2	Methylene chloride	ND	570
91-20-3	Naphthalene	ND	570
103-65-1	n-Propylbenzene	ND	570
100-42-5	Styrene	ND	570
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	57
109-99-9	Tetrahydrofuran	ND	570
96-18-4	1,2,3-Trichloropropane	ND	570
630-20-6	1,1,1,2-Tetrachloroethane	ND	570
79-34-5	1,1,2,2-Tetrachloroethane	ND	570
127-18-4	Tetrachloroethene (PCE)	15000	570
108-88-3	Toluene	ND	570
87-61-6	1,2,3-Trichlorobenzene	ND	570
120-82-1	1,2,4-Trichlorobenzene	ND	570
71-55-6	1,1,1-Trichloroethane	ND	570
79-00-5	1,1,2-Trichloroethane	ND	570
79-01-6	Trichloroethene (TCE)	19000	570
75-69-4	Trichlorofluoromethane	ND	1100
95-63-6	1,2,4-Trimethylbenzene	ND	570
108-67-8	1,3,5-Trimethylbenzene	ND	570
75-01-4	Vinyl chloride	ND	1100
95-47-6	o-Xylene	ND	570
	m,p-Xylenes	ND	570
Surrogate	Recovery	Limits	
Bromofluorobenzene	90%	78%-111%	
1,2-Dichloroethane-d4	99%	91%-114%	
TP 1 10	10.407	0.60/ 11.50/	

Surrogate	Recovery	Limits
Bromofluorobenzene	90%	78%-111%
1,2-Dichloroethane-d4	99%	91%-114%
Toluene-d8	104%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	8	Sample Description:	841061130-08
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	11.7
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6900	ug/kg
C9-C12 Aliphatics**	50	ND	6900	ug/kg
C9-C10 Aromatics***	50	ND	6900	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	95	70%-130%
2,5-dibromotoluene #2	99	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
Benzene	ND	340	ug/kg
Ethylbenzene	ND	340	ug/kg
Methyl tert-butyl ether (MTBE)	ND	69	ug/kg
Naphthalene	ND	340	ug/kg
Toluene	ND	340	ug/kg
m,p-Xylenes	ND	340	ug/kg
o-Xylene	ND	340	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 Sample Description: 841061130-08

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:11.7Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/07/06By:Dilution Factor:200

Method: 8260B Soil Extract Volume:
QC Batch#: 51101 Lab Data File: J28611.D;M32621.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	4500
71-43-2	Benzene	ND	1100
108-86-1	Bromobenzene	ND	1100
74-97-5	Bromochloromethane	ND	1100
75-27-4	Bromodichloromethane	ND	1100
75-25-2	Bromoform	ND	1100
74-83-9	Bromomethane	ND	2300
78-93-3	2-Butanone (MEK)	ND	2300
104-51-8	n-Butylbenzene	ND	1100
135-98-8	sec-Butylbenzene	ND	1100
98-06-6	tert-Butylbenzene	ND	1100
75-15-0	Carbon disulfide	ND	1100
56-23-5	Carbon tetrachloride	ND	1100
108-90-7	Chlorobenzene	ND	1100
75-00-3	Chloroethane	ND	2300
67-66-3	Chloroform	ND	1100
74-87-3	Chloromethane	ND	2300
95-49-8	2-Chlorotoluene	ND	1100
106-43-4	4-Chlorotoluene	ND	1100
108-20-3	Di-isopropyl ether (DIPE)	ND	57
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	1100
124-48-1	Dibromochloromethane	ND	1100
106-93-4	1,2-Dibromoethane (EDB)	ND	680
74-95-3	Dibromomethane	ND	1100
95-50-1	1,2-Dichlorobenzene	ND	1100
541-73-1	1,3-Dichlorobenzene	ND	1100
106-46-7	1,4-Dichlorobenzene	ND	1100
75-71-8	Dichlorodifluoromethane	ND	2300
75-34-3	1,1-Dichloroethane	ND	1100
107-06-2	1,2-Dichloroethane	ND	1100
75-35-4	1,1-Dichloroethene	ND	1100
156-59-2	cis-1,2-Dichloroethene	ND	1100
156-60-5	trans-1,2-Dichloroethene	ND	1100
78-87-5	1,2-Dichloropropane	ND	1100
142-28-9	1,3-Dichloropropane	ND	1100
590-20-7	2,2-Dichloropropane	ND	1100
563-58-6	1,1-Dichloropropene	ND	1100
10061-01-5	cis-1,3-Dichloropropene	ND	1100
10061-02-6	trans-1,3-Dichloropropene	ND	1100
60-29-7	Diethyl ether	ND	2300
123-91-1	1,4-Dioxane	ND	4500

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 (continued) Sample Description: 841061130-08

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 11.7

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/07/06 By: GP Dilution Factor: 200

Method: 8260B Soil Extract Volume:

Method: 8260B Soil Extract Volume:
QC Batch#: 51101 Lab Data File: J28611.D;M32621.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	57
100-41-4	Ethylbenzene	ND	1100
87-68-3	Hexachlorobutadiene	ND	1100
591-78-6	2-Hexanone	ND	2300
98-82-8	Isopropylbenzene	ND	1100
99-87-6	4-Isopropyltoluene	ND	1100
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1100
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	2300
75-09-2	Methylene chloride	ND	1100
91-20-3	Naphthalene	ND	1100
103-65-1	n-Propylbenzene	ND	1100
100-42-5	Styrene	ND	1100
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	57
109-99-9	Tetrahydrofuran	ND	1100
96-18-4	1,2,3-Trichloropropane	ND	1100
630-20-6	1,1,1,2-Tetrachloroethane	ND	1100
79-34-5	1,1,2,2-Tetrachloroethane	ND	1100
127-18-4	Tetrachloroethene (PCE)	20000	1100
108-88-3	Toluene	ND	1100
87-61-6	1,2,3-Trichlorobenzene	ND	1100
120-82-1	1,2,4-Trichlorobenzene	ND	1100
71-55-6	1,1,1-Trichloroethane	ND	1100
79-00-5	1,1,2-Trichloroethane	ND	1100
79-01-6	Trichloroethene (TCE)	31000	1100
75-69-4	Trichlorofluoromethane	ND	2300
95-63-6	1,2,4-Trimethylbenzene	ND	1100
108-67-8	1,3,5-Trimethylbenzene	ND	1100
75-01-4	Vinyl chloride	ND	2300
95-47-6	o-Xylene	ND	1100
	m,p-Xylenes	ND	1100
Surrogate	Recovery	Limits	
Bromofluorobenzene	88%	87%-105%	
1,2-Dichloroethane-d4	102%	91%-109%	
Toluene-d8	97%	92%-105%	

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	9	Sample Description:	841061130-09
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	5.6
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	5900	ug/kg
C9-C12 Aliphatics**	50	ND	5900	ug/kg
C9-C10 Aromatics***	50	ND	5900	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	93	70%-130%	
2,5-dibromotoluene #2	98	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	290	ug/kg
Ethylbenzene	ND	290	ug/kg
Methyl tert-butyl ether (MTBE)	ND	59	ug/kg
Naphthalene	ND	290	ug/kg
Toluene	ND	290	ug/kg
m,p-Xylenes	ND	290	ug/kg
o-Xylene	ND	290	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 9 Sample Description: 841061130-09

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:5.6Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28536.D;M32622.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	20
71-43-2	Benzene	ND	5.0
108-86-1	Bromobenzene	ND	5.0
74-97-5	Bromochloromethane	ND	5.0
75-27-4	Bromodichloromethane	ND	5.0
75-25-2	Bromoform	ND	5.0
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.0
135-98-8	sec-Butylbenzene	ND	5.0
98-06-6	tert-Butylbenzene	ND	5.0
75-15-0	Carbon disulfide	ND	5.0
56-23-5	Carbon tetrachloride	ND	5.0
108-90-7	Chlorobenzene	ND	5.0
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.0
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.0
106-43-4	4-Chlorotoluene	ND	5.0
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.0
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.0
74-95-3	Dibromomethane	ND	5.0
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.0
106-46-7	1,4-Dichlorobenzene	ND	5.0
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.0
107-06-2	1,2-Dichloroethane	ND	5.0
75-35-4	1,1-Dichloroethene	ND	5.0
156-59-2	cis-1,2-Dichloroethene	ND	5.0
156-60-5	trans-1,2-Dichloroethene	ND	5.0
78-87-5	1,2-Dichloropropane	ND	5.0
142-28-9	1,3-Dichloropropane	ND	5.0
590-20-7	2,2-Dichloropropane	ND	5.0
563-58-6	1,1-Dichloropropene	ND	5.0
10061-01-5	cis-1,3-Dichloropropene	ND	5.0
10061-02-6	trans-1,3-Dichloropropene	ND	5.0
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 9 (continued) Sample Description: 841061130-09

Date Collected: 11/30/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 5.6 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/05/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume:

Lab Data File: J28536.D;M32622.D QC Batch#: 50075

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	50
100-41-4	Ethylbenzene	ND	5.0
87-68-3	Hexachlorobutadiene	ND	5.0
591-78-6	2-Hexanone	ND	10
98-82-8	Isopropylbenzene	ND	5.0
99-87-6	4-Isopropyltoluene	ND	5.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	5.0
103-65-1	n-Propylbenzene	ND	5.0
100-42-5	Styrene	ND	5.0
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	50
109-99-9	Tetrahydrofuran	ND	5.0
96-18-4	1,2,3-Trichloropropane	ND	5.0
630-20-6	1,1,2-Tetrachloroethane	ND	5.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	110	5.0
108-88-3	Toluene	ND	5.0
87-61-6	1,2,3-Trichlorobenzene	ND	5.0
120-82-1	1,2,4-Trichlorobenzene	ND	5.0
71-55-6	1,1,1-Trichloroethane	ND	5.0
79-00-5	1,1,2-Trichloroethane	ND	5.0
79-01-6	Trichloroethene (TCE)	58	5.0
75-69-4	Trichlorofluoromethane	ND	10
95-63-6	1,2,4-Trimethylbenzene	ND	5.0
108-67-8	1,3,5-Trimethylbenzene	ND	5.0
75-01-4	Vinyl chloride	ND	10
95-47-6	o-Xylene	ND	5.0
	m,p-Xylenes	ND	5.0
Surrogate	Recovery	Limits	
Bromofluorobenzene	78%	78%-111%	
1,2-Dichloroethane-d4	109%	91%-114%	
TD 1 10	1070/	0.60/ 11.50/	

Surrogate	Recovery	Limits
Bromofluorobenzene	78%	78%-111%
1,2-Dichloroethane-d4	109%	91%-114%
Toluene-d8	107%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	10	Sample Description:	841061130-10
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	8.1
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6500	ug/kg
C9-C12 Aliphatics**	50	ND	6500	ug/kg
C9-C10 Aromatics***	50	ND	6500	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	99	70%-130%	
2,5-dibromotoluene #2	105	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	330	ug/kg
Ethylbenzene	ND	330	ug/kg
Methyl tert-butyl ether (MTBE)	ND	65	ug/kg
Naphthalene	ND	330	ug/kg
Toluene	ND	330	ug/kg
m,p-Xylenes	ND	330	ug/kg
o-Xylene	ND	330	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 10 Sample Description: 841061130-10

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 8.1

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50075 Lab Data File: J28537.D;M32623.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	21
71-43-2	Benzene	ND	5.3
108-86-1	Bromobenzene	ND	5.3
74-97-5	Bromochloromethane	ND	5.3
75-27-4	Bromodichloromethane	ND	5.3
75-25-2	Bromoform	ND	5.3
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.3
135-98-8	sec-Butylbenzene	ND	5.3
98-06-6	tert-Butylbenzene	ND	5.3
75-15-0	Carbon disulfide	ND	5.3
56-23-5	Carbon tetrachloride	ND	5.3
108-90-7	Chlorobenzene	ND	5.3
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.3
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.3
106-43-4	4-Chlorotoluene	ND	5.3
108-20-3	Di-isopropyl ether (DIPE)	ND	53
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.3
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
74-95-3	Dibromomethane	ND	5.3
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.3
106-46-7	1,4-Dichlorobenzene	ND	5.3
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.3
107-06-2	1,2-Dichloroethane	ND	5.3
75-35-4	1,1-Dichloroethene	ND	5.3
156-59-2	cis-1,2-Dichloroethene	ND	5.3
156-60-5	trans-1,2-Dichloroethene	ND	5.3
78-87-5	1,2-Dichloropropane	ND	5.3
142-28-9	1,3-Dichloropropane	ND	5.3
590-20-7	2,2-Dichloropropane	ND	5.3
563-58-6	1,1-Dichloropropene	ND	5.3
10061-01-5	cis-1,3-Dichloropropene	ND	5.3
10061-02-6	trans-1,3-Dichloropropene	ND	5.3
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	21

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 10 (continued) Sample Description: 841061130-10

Date Collected: 11/30/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 8.1 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/05/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume:

QC Batch#: 50075 Lab Data File: J28537.D;M32623.D

Units: ug/kg

CAS No.	AS No. Parameter		esult	DL
	Ethyl tertiary-butyl ether (EtBE)		ND	53
100-41-4	Ethylbenzene		ND	5.3
87-68-3	Hexachlorobutadiene		ND	5.3
591-78-6	2-Hexanone		ND	10
98-82-8	Isopropylbenzene		ND	5.3
99-87-6	4-Isopropyltoluene		ND	5.3
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	5.3
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	10
75-09-2	Methylene chloride		ND	5.3
91-20-3	Naphthalene		ND	5.3
103-65-1	n-Propylbenzene		ND	5.3
100-42-5	Styrene		ND	5.3
994-05-8	Tertiary-amyl methyl ether (TAME)		ND	53
109-99-9	Tetrahydrofuran		ND	5.3
96-18-4	1,2,3-Trichloropropane		ND	5.3
630-20-6	1,1,1,2-Tetrachloroethane		ND	5.3
79-34-5	1,1,2,2-Tetrachloroethane		ND	5.0
127-18-4	Tetrachloroethene (PCE)		22	5.3
108-88-3	Toluene		ND	5.3
87-61-6	1,2,3-Trichlorobenzene		ND	5.3
120-82-1	1,2,4-Trichlorobenzene		ND	5.3
71-55-6	1,1,1-Trichloroethane		ND	5.3
79-00-5	1,1,2-Trichloroethane		ND	5.3
79-01-6	Trichloroethene (TCE)		9.6	5.3
75-69-4	Trichlorofluoromethane		ND	10
95-63-6	1,2,4-Trimethylbenzene		ND	5.3
108-67-8	1,3,5-Trimethylbenzene		ND	5.3
75-01-4	Vinyl chloride		ND	10
95-47-6	o-Xylene		ND	5.3
	m,p-Xylenes		ND	5.3
Surrogate	Recovery	Limits		
Bromofluorobenzene	93%	78%-111%		
1,2-Dichloroethane-d4	105%	91%-114%		
Toluono de	050/	960/ 1150/		

Surrogate	Recovery	Limits
Bromofluorobenzene	93%	78%-111%
1,2-Dichloroethane-d4	105%	91%-114%
Toluene-d8	95%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	11	Sample Description:	841061130-11
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	6.6
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	5900	ug/kg
C9-C12 Aliphatics**	50	ND	5900	ug/kg
C9-C10 Aromatics***	50	ND	5900	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	92	70%-130%
2,5-dibromotoluene #2	100	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	300	ug/kg
Ethylbenzene	ND	300	ug/kg
Methyl tert-butyl ether (MTBE)	ND	59	ug/kg
Naphthalene	ND	300	ug/kg
Toluene	ND	300	ug/kg
m,p-Xylenes	ND	300	ug/kg
o-Xylene	ND	300	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 11 Sample Description: 841061130-11

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 6.6

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/07/06 By: GP Dilution Factor: 200

Method: 8260B Soil Extract Volume:

QC Batch#: 51101 Lab Data File: J28600.D;M32624.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	4300
71-43-2	Benzene	ND	1100
108-86-1	Bromobenzene	ND	1100
74-97-5	Bromochloromethane	ND	1100
75-27-4	Bromodichloromethane	ND	1100
75-25-2	Bromoform	ND	1100
74-83-9	Bromomethane	ND	2100
78-93-3	2-Butanone (MEK)	ND	2100
104-51-8	n-Butylbenzene	ND	1100
135-98-8	sec-Butylbenzene	ND	1100
98-06-6	tert-Butylbenzene	ND	1100
75-15-0	Carbon disulfide	ND	1100
56-23-5	Carbon tetrachloride	ND	1100
108-90-7	Chlorobenzene	ND	1100
75-00-3	Chloroethane	ND	2100
67-66-3	Chloroform	ND	1100
74-87-3	Chloromethane	ND	2100
95-49-8	2-Chlorotoluene	ND	1100
106-43-4	4-Chlorotoluene	ND	1100
108-20-3	Di-isopropyl ether (DIPE)	ND	54
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	1100
124-48-1	Dibromochloromethane	ND	1100
106-93-4	1,2-Dibromoethane (EDB)	ND	640
74-95-3	Dibromomethane	ND	1100
95-50-1	1,2-Dichlorobenzene	ND	1100
541-73-1	1,3-Dichlorobenzene	ND	1100
106-46-7	1,4-Dichlorobenzene	ND	1100
75-71-8	Dichlorodifluoromethane	ND	2100
75-34-3	1,1-Dichloroethane	ND	1100
107-06-2	1,2-Dichloroethane	ND	1100
75-35-4	1,1-Dichloroethene	ND	1100
156-59-2	cis-1,2-Dichloroethene	ND	1100
156-60-5	trans-1,2-Dichloroethene	ND	1100
78-87-5	1,2-Dichloropropane	ND	1100
142-28-9	1,3-Dichloropropane	ND	1100
590-20-7	2,2-Dichloropropane	ND	1100
563-58-6	1,1-Dichloropropene	ND	1100
10061-01-5	cis-1,3-Dichloropropene	ND	1100
10061-02-6	trans-1,3-Dichloropropene	ND	1100
60-29-7	Diethyl ether	ND	2100
123-91-1	1,4-Dioxane	ND	4300

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 11 (continued) Sample Description: 841061130-11

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:6.6Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/07/06By:GPDilution Factor:200Method:8260BSoil Extract Volume:

QC Batch#: 51101 Lab Data File: J28600.D;M32624.D

Units: ug/kg

956D450

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	54
100-41-4	Ethylbenzene	ND	1100
87-68-3	Hexachlorobutadiene	ND	1100
591-78-6	2-Hexanone	ND	2100
98-82-8	Isopropylbenzene	ND	1100
99-87-6	4-Isopropyltoluene	ND	1100
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1100
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	2100
75-09-2	Methylene chloride	ND	1100
91-20-3	Naphthalene	ND	1100
103-65-1	n-Propylbenzene	ND	1100
100-42-5	Styrene	ND	1100
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	54
109-99-9	Tetrahydrofuran	ND	1100
96-18-4	1,2,3-Trichloropropane	ND	1100
630-20-6	1,1,1,2-Tetrachloroethane	ND	1100
79-34-5	1,1,2,2-Tetrachloroethane	ND	1100
127-18-4	Tetrachloroethene (PCE)	34000	1100
108-88-3	Toluene	ND	1100
87-61-6	1,2,3-Trichlorobenzene	ND	1100
120-82-1	1,2,4-Trichlorobenzene	ND	1100
71-55-6	1,1,1-Trichloroethane	ND	1100
79-00-5	1,1,2-Trichloroethane	ND	1100
79-01-6	Trichloroethene (TCE)	6700	1100
75-69-4	Trichlorofluoromethane	ND	2100
95-63-6	1,2,4-Trimethylbenzene	ND	1100
108-67-8	1,3,5-Trimethylbenzene	ND	1100
75-01-4	Vinyl chloride	ND	2100
95-47-6	o-Xylene	ND	1100
	m,p-Xylenes	ND	1100
Surrogate	Recovery	Limits	
Bromofluorobenzene	89%	87%-105%	
1,2-Dichloroethane-d4	101%	91%-109%	
Toluene-d8	98%	92%-105%	

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	12	Sample Description:	841061130-12
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	14.7
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7800	ug/kg
C9-C12 Aliphatics**	50	ND	7800	ug/kg
C9-C10 Aromatics***	50	ND	7800	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	99	70%-130%
2,5-dibromotoluene #2	107	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	390	ug/kg
Ethylbenzene	ND	390	ug/kg
Methyl tert-butyl ether (MTBE)	ND	78	ug/kg
Naphthalene	ND	390	ug/kg
Toluene	ND	390	ug/kg
m,p-Xylenes	ND	390	ug/kg
o-Xylene	ND	390	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 12 Sample Description: 841061130-12

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:14.7Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28538.D;M32625.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	23
71-43-2	Benzene	ND	5.8
108-86-1	Bromobenzene	ND	5.8
74-97-5	Bromochloromethane	ND	5.8
75-27-4	Bromodichloromethane	ND	5.8
75-25-2	Bromoform	ND	5.8
74-83-9	Bromomethane	ND	12
78-93-3	2-Butanone (MEK)	ND	12
104-51-8	n-Butylbenzene	ND	5.8
135-98-8	sec-Butylbenzene	ND	5.8
98-06-6	tert-Butylbenzene	ND	5.8
75-15-0	Carbon disulfide	ND	5.8
56-23-5	Carbon tetrachloride	ND	5.8
108-90-7	Chlorobenzene	ND	5.8
75-00-3	Chloroethane	ND	12
67-66-3	Chloroform	ND	5.8
74-87-3	Chloromethane	ND	12
95-49-8	2-Chlorotoluene	ND	5.8
106-43-4	4-Chlorotoluene	ND	5.8
108-20-3	Di-isopropyl ether (DIPE)	ND	58
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.8
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.5
74-95-3	Dibromomethane	ND	5.8
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.8
106-46-7	1,4-Dichlorobenzene	ND	5.8
75-71-8	Dichlorodifluoromethane	ND	12
75-34-3	1,1-Dichloroethane	ND	5.8
107-06-2	1,2-Dichloroethane	ND	5.8
75-35-4	1,1-Dichloroethene	ND	5.8
156-59-2	cis-1,2-Dichloroethene	ND	5.8
156-60-5	trans-1,2-Dichloroethene	ND	5.8
78-87-5	1,2-Dichloropropane	ND	5.8
142-28-9	1,3-Dichloropropane	ND	5.8
590-20-7	2,2-Dichloropropane	ND	5.8
563-58-6	1,1-Dichloropropene	ND	5.8
10061-01-5	cis-1,3-Dichloropropene	ND	5.8
10061-02-6	trans-1,3-Dichloropropene	ND	5.8
60-29-7	Diethyl ether	ND	12
123-91-1	1,4-Dioxane	ND	23
	,		-

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 12 (continued) Sample Description: 841061130-12

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 14.7

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50075 Lab Data File: J28538.D;M32625.D

Units: ug/kg

CAS No.	S No. Parameter		DL
	Ethyl tertiary-butyl ether (EtBE)	ND	58
100-41-4	Ethylbenzene	ND	5.8
87-68-3	Hexachlorobutadiene	ND	5.8
591-78-6	2-Hexanone	ND	12
98-82-8	Isopropylbenzene	ND	5.8
99-87-6	4-Isopropyltoluene	ND	5.8
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.8
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	12
75-09-2	Methylene chloride	ND	5.8
91-20-3	Naphthalene	ND	5.8
103-65-1	n-Propylbenzene	ND	5.8
100-42-5	Styrene	ND	5.8
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	58
109-99-9	Tetrahydrofuran	ND	5.8
96-18-4	1,2,3-Trichloropropane	ND	5.8
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.8
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	310	5.8
108-88-3	Toluene	ND	5.8
87-61-6	1,2,3-Trichlorobenzene	ND	5.8
120-82-1	1,2,4-Trichlorobenzene	ND	5.8
71-55-6	1,1,1-Trichloroethane	ND	5.8
79-00-5	1,1,2-Trichloroethane	ND	5.8
79-01-6	Trichloroethene (TCE)	79	5.8
75-69-4	Trichlorofluoromethane	ND	12
95-63-6	1,2,4-Trimethylbenzene	ND	5.8
108-67-8	1,3,5-Trimethylbenzene	ND	5.8
75-01-4	Vinyl chloride	ND	12
95-47-6	o-Xylene	ND	5.8
	m,p-Xylenes	ND	5.8
Surrogate	Recovery Lir	nits	
Bromofluorobenzene	87% 78	%-111%	

Surrogate	Recovery	Limits
Bromofluorobenzene	87%	78%-111%
1,2-Dichloroethane-d4	100%	91%-114%
Toluene-d8	101%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	13	Sample Description:	841061130-13
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	8.6
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6400	ug/kg
C9-C12 Aliphatics**	50	ND	6400	ug/kg
C9-C10 Aromatics***	50	ND	6400	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	97	70%-130%
2,5-dibromotoluene #2	107	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	320	ug/kg
Ethylbenzene	ND	320	ug/kg
Methyl tert-butyl ether (MTBE)	ND	64	ug/kg
Naphthalene	ND	320	ug/kg
Toluene	ND	320	ug/kg
m,p-Xylenes	ND	320	ug/kg
o-Xylene	ND	320	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Soil Extract Volume:

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 13 Sample Description: 841061130-13

Date Collected: 11/30/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 8.6 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/06/06 By: GP Dilution Factor: 1 Method: 8260B

QC Batch#: 50999 Lab Data File: J28553.D;M32626.D

Units: ug/kg

CAS No. Parameter	Result	DL
67-64-1 Acetone	ND	21
71-43-2 Benzene	ND	5.2
108-86-1 Bromobenzene	ND	5.2
74-97-5 Bromochloromethan	ne ND	5.2
75-27-4 Bromodichlorometh	nane ND	5.2
75-25-2 Bromoform	ND	5.2
74-83-9 Bromomethane	ND	10
78-93-3 2-Butanone (MEK)	ND	10
104-51-8 n-Butylbenzene	ND	5.2
135-98-8 sec-Butylbenzene	ND	5.2
98-06-6 tert-Butylbenzene	ND	5.2
75-15-0 Carbon disulfide	ND	5.2
56-23-5 Carbon tetrachloride	e ND	5.2
108-90-7 Chlorobenzene	ND	5.2
75-00-3 Chloroethane	ND	10
67-66-3 Chloroform	ND	5.2
74-87-3 Chloromethane	ND	10
95-49-8 2-Chlorotoluene	ND	5.2
106-43-4 4-Chlorotoluene	ND	5.2
108-20-3 Di-isopropyl ether (I	DIPE) ND	52
96-12-8 1,2-Dibromo-3-chlo		5.2
124-48-1 Dibromochlorometh	nane ND	5.0
106-93-4 1,2-Dibromoethane ((EDB) ND	3.1
74-95-3 Dibromomethane	ND	5.2
95-50-1 1,2-Dichlorobenzen	e ND	5.0
541-73-1 1,3-Dichlorobenzen	e ND	5.2
106-46-7 1,4-Dichlorobenzen	e ND	5.2
75-71-8 Dichlorodifluorome	ethane ND	10
75-34-3 1,1-Dichloroethane	ND	5.2
107-06-2 1,2-Dichloroethane	ND	5.2
75-35-4 1,1-Dichloroethene	ND	5.2
156-59-2 cis-1,2-Dichloroethe	ene ND	5.2
156-60-5 trans-1,2-Dichloroe	thene ND	5.2
78-87-5 1,2-Dichloropropan	e ND	5.2
142-28-9 1,3-Dichloropropan		5.2
590-20-7 2,2-Dichloropropan	e ND	5.2
563-58-6 1,1-Dichloropropen		5.2
10061-01-5 cis-1,3-Dichloropro		5.2
10061-02-6 trans-1,3-Dichlorop	•	5.2
60-29-7 Diethyl ether	ND	10
	1,2	10

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E612052 PL Sample No: 13 (continued) Sample Description: 841061130-13

Date Collected: 11/30/2006 Matrix: Solid Date Received: 12/1/2006 Percent Moisture: 8.6 Sample Weight/Volume: Date Extracted: By: Date Analyzed: 12/06/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume:

QC Batch#: 50999 Lab Data File: J28553.D;M32626.D

Units: ug/kg

CAS No.	Parameter		DL
	Ethyl tertiary-butyl ether (EtBE)	ND	52
100-41-4	Ethylbenzene	ND	5.2
87-68-3	Hexachlorobutadiene	ND	5.2
591-78-6	2-Hexanone	ND	10
98-82-8	Isopropylbenzene	ND	5.2
99-87-6	4-Isopropyltoluene	ND	5.2
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.2
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10
75-09-2	Methylene chloride	ND	5.2
91-20-3	Naphthalene	ND	5.2
103-65-1	n-Propylbenzene	ND	5.2
100-42-5	Styrene	ND	5.2
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	52
109-99-9	Tetrahydrofuran	ND	5.2
96-18-4	1,2,3-Trichloropropane	ND	5.2
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.2
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	130	5.2
108-88-3	Toluene	Toluene ND	
87-61-6	1,2,3-Trichlorobenzene		
120-82-1	1,2,4-Trichlorobenzene	1,2,4-Trichlorobenzene ND	
71-55-6	1,1,1-Trichloroethane	1,1,1-Trichloroethane ND	
79-00-5	1,1,2-Trichloroethane	ND	5.2
79-01-6	Trichloroethene (TCE)	150	5.2
75-69-4	Trichlorofluoromethane	ND	10
95-63-6	1,2,4-Trimethylbenzene	ND	5.2
108-67-8	1,3,5-Trimethylbenzene	ND	5.2
75-01-4	Vinyl chloride	ND	10
95-47-6	o-Xylene	ND	5.2
	m,p-Xylenes	ND	5.2
Surrogate	Recovery	Limits	
Bromofluorobenzene	66%	78%-111%	
1,2-Dichloroethane-d4	104%	91%-114%	

Surrogate	Recovery	Limits
Bromofluorobenzene	66%	78%-111%
1,2-Dichloroethane-d4	104%	91%-114%
Toluene-d8	117%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	14	Sample Description:	841061130-14
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	8.9
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6400	ug/kg
C9-C12 Aliphatics**	50	ND	6400	ug/kg
C9-C10 Aromatics***	50	ND	6400	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	98	70%-130%
2,5-dibromotoluene #2	105	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	320	ug/kg
Ethylbenzene	ND	320	ug/kg
Methyl tert-butyl ether (MTBE)	ND	64	ug/kg
Naphthalene	ND	320	ug/kg
Toluene	ND	320	ug/kg
m,p-Xylenes	ND	320	ug/kg
o-Xylene	ND	320	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 14 Sample Description: 841061130-14

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:8.9Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28540.D;J28554.;M32627.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	22
71-43-2	Benzene	ND	5.4
108-86-1	Bromobenzene	ND	5.4
74-97-5	Bromochloromethane	ND	5.4
75-27-4	Bromodichloromethane	ND	5.4
75-25-2	Bromoform	ND	5.4
74-83-9	Bromomethane	ND	11
78-93-3	2-Butanone (MEK)	ND	11
104-51-8	n-Butylbenzene	ND	5.4
135-98-8	sec-Butylbenzene	ND	5.4
98-06-6	tert-Butylbenzene	ND	5.4
75-15-0	Carbon disulfide	ND	5.4
56-23-5	Carbon tetrachloride	ND	5.4
108-90-7	Chlorobenzene	ND	5.4
75-00-3	Chloroethane	ND	11
67-66-3	Chloroform	ND	5.4
74-87-3	Chloromethane	ND	11
95-49-8	2-Chlorotoluene	ND	5.4
106-43-4	4-Chlorotoluene	ND	5.4
108-20-3	Di-isopropyl ether (DIPE)	ND	54
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.4
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
74-95-3	Dibromomethane	ND	5.4
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.4
106-46-7	1,4-Dichlorobenzene	ND	5.4
75-71-8	Dichlorodifluoromethane	ND	11
75-34-3	1,1-Dichloroethane	ND	5.4
107-06-2	1,2-Dichloroethane	ND	5.4
75-35-4	1,1-Dichloroethene	ND	5.4
156-59-2	cis-1,2-Dichloroethene	ND	5.4
156-60-5	trans-1,2-Dichloroethene	ND	5.4
78-87-5	1,2-Dichloropropane	ND	5.4
142-28-9	1,3-Dichloropropane	ND	5.4
590-20-7	2,2-Dichloropropane	ND	5.4
563-58-6	1,1-Dichloropropene	ND	5.4
10061-01-5	cis-1,3-Dichloropropene	ND	5.4
10061-02-6	trans-1,3-Dichloropropene	ND	5.4
60-29-7	Diethyl ether	ND	11
123-91-1	1,4-Dioxane	ND	22

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 14 (continued) Sample Description: 841061130-14

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:8.9Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/05/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50075 Lab Data File: J28540.D;J28554.;M32627.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	54
100-41-4	Ethylbenzene	ND	5.4
87-68-3	Hexachlorobutadiene	ND	5.4
591-78-6	2-Hexanone	ND	11
98-82-8	Isopropylbenzene	ND	5.4
99-87-6	4-Isopropyltoluene	ND	5.4
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.4
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	11
75-09-2	Methylene chloride	ND	5.4
91-20-3	Naphthalene	ND	5.4
103-65-1	n-Propylbenzene	ND	5.4
100-42-5	Styrene	ND	5.4
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	54
109-99-9	Tetrahydrofuran	ND	5.4
96-18-4	1,2,3-Trichloropropane	ND	5.4
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.4
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	120	5.4
108-88-3	Toluene	ND	5.4
87-61-6	1,2,3-Trichlorobenzene	ND	5.4
120-82-1	1,2,4-Trichlorobenzene	ND	5.4
71-55-6	1,1,1-Trichloroethane	ND	5.4
79-00-5	1,1,2-Trichloroethane	ND	5.4
79-01-6	Trichloroethene (TCE)	67	5.4
75-69-4	Trichlorofluoromethane	ND	11
95-63-6	1,2,4-Trimethylbenzene	ND	5.4
108-67-8	1,3,5-Trimethylbenzene	ND	5.4
75-01-4	Vinyl chloride	ND	11
95-47-6	o-Xylene	ND	5.4
	m,p-Xylenes	ND	5.4
Surrogate	Recovery Lir	nits	
Bromofluorobenzene	79% 78	%-111%	

Surrogate	Recovery	Limits
Bromofluorobenzene	79%	78%-111%
1,2-Dichloroethane-d4	106%	91%-114%
Toluene-d8	104%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	15	Sample Description:	841061130-15
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	11.6
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	7600	ug/kg
C9-C12 Aliphatics**	50	ND	7600	ug/kg
C9-C10 Aromatics***	50	ND	7600	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	97	70%-130%
2,5-dibromotoluene #2	105	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	380	ug/kg
Ethylbenzene	ND	380	ug/kg
Methyl tert-butyl ether (MTBE)	ND	76	ug/kg
Naphthalene	ND	380	ug/kg
Toluene	ND	380	ug/kg
m,p-Xylenes	ND	380	ug/kg
o-Xylene	ND	380	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Matrix: Solid

Percent Moisture: 11.6 Sample Weight/Volume:

Dilution Factor: 1

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 15 Sample Description: 841061130-15

Date Collected: 11/30/2006

Date Received: 12/1/2006

Date Extracted: By:
Date Analyzed: 12/05/06 By: GP

Method: 8260B Soil Extract Volume:
QC Batch#: 50075 Lab Data File: J28541.D;M32628.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	21
71-43-2	Benzene	ND	5.2
108-86-1	Bromobenzene	ND	5.2
74-97-5	Bromochloromethane	ND	5.2
75-27-4	Bromodichloromethane	ND	5.2
75-25-2	Bromoform	ND	5.2
74-83-9	Bromomethane	ND	10
78-93-3	2-Butanone (MEK)	ND	10
104-51-8	n-Butylbenzene	ND	5.2
135-98-8	sec-Butylbenzene	ND	5.2
98-06-6	tert-Butylbenzene	ND	5.2
75-15-0	Carbon disulfide	ND	5.2
56-23-5	Carbon tetrachloride	ND	5.2
108-90-7	Chlorobenzene	ND	5.2
75-00-3	Chloroethane	ND	10
67-66-3	Chloroform	ND	5.2
74-87-3	Chloromethane	ND	10
95-49-8	2-Chlorotoluene	ND	5.2
106-43-4	4-Chlorotoluene	ND	5.2
108-20-3	Di-isopropyl ether (DIPE)	ND	52
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.2
124-48-1	Dibromochloromethane	ND	5.0
106-93-4	1,2-Dibromoethane (EDB)	ND	3.1
74-95-3	Dibromomethane	ND	5.2
95-50-1	1,2-Dichlorobenzene	ND	5.0
541-73-1	1,3-Dichlorobenzene	ND	5.2
106-46-7	1,4-Dichlorobenzene	ND	5.2
75-71-8	Dichlorodifluoromethane	ND	10
75-34-3	1,1-Dichloroethane	ND	5.2
107-06-2	1,2-Dichloroethane	ND	5.2
75-35-4	1,1-Dichloroethene	ND	5.2
156-59-2	cis-1,2-Dichloroethene	ND	5.2
156-60-5	trans-1,2-Dichloroethene	ND	5.2
78-87-5	1,2-Dichloropropane	ND	5.2
142-28-9	1,3-Dichloropropane	ND	5.2
590-20-7	2,2-Dichloropropane	ND	5.2
563-58-6	1,1-Dichloropropene	ND	5.2
10061-01-5	cis-1,3-Dichloropropene	ND	5.2
10061-02-6	trans-1,3-Dichloropropene	ND	5.2
60-29-7	Diethyl ether	ND	10
123-91-1	1,4-Dioxane	ND	21

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Matrix: Solid

Percent Moisture: 11.6 Sample Weight/Volume:

Customer: Fuss & O'Neill Laboratory: Premier Laboratory, LLC Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 15 (continued) Sample Description: 841061130-15

Date Collected: 11/30/2006 Date Received: 12/1/2006 Date Extracted: By:

Date Analyzed: 12/05/06 By: GP Dilution Factor: 1 Method: 8260B Soil Extract Volume: QC Batch#: 50075 Lab Data File: J28541.D;M32628.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	52
100-41-4	Ethylbenzene	ND	5.2
87-68-3	Hexachlorobutadiene	ND	5.2
591-78-6	2-Hexanone	ND	10
98-82-8	Isopropylbenzene	ND	5.2
99-87-6	4-Isopropyltoluene	ND	5.2
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.2
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	10
75-09-2	Methylene chloride	ND	5.2
91-20-3	Naphthalene	ND	5.2
103-65-1	n-Propylbenzene	ND	5.2
100-42-5	Styrene	ND	5.2
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	52
109-99-9	Tetrahydrofuran	ND	5.2
96-18-4	1,2,3-Trichloropropane	ND	5.2
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.2
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.0
127-18-4	Tetrachloroethene (PCE)	13	5.2
108-88-3	Toluene	ND	5.2
87-61-6	1,2,3-Trichlorobenzene	ND	5.2
120-82-1	1,2,4-Trichlorobenzene	ND	5.2
71-55-6	1,1,1-Trichloroethane	1,1,1-Trichloroethane 73	
79-00-5	1,1,2-Trichloroethane	ND	5.2
79-01-6	Trichloroethene (TCE)	37	5.2
75-69-4	Trichlorofluoromethane	ND	10
95-63-6	1,2,4-Trimethylbenzene	ND	5.2
108-67-8	1,3,5-Trimethylbenzene	ND	5.2
75-01-4	Vinyl chloride	ND	10
95-47-6	o-Xylene	ND	5.2
	m,p-Xylenes	ND	5.2
Surrogate	Recovery	Limits	
Bromofluorobenzene	43%	78%-111%	
1,2-Dichloroethane-d4	105%	91%-114%	

Toluene-d8 148% 86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	16	Sample Description:	841061130-16
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	6.2
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6000	ug/kg
C9-C12 Aliphatics**	50	ND	6000	ug/kg
C9-C10 Aromatics***	50	ND	6000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	109	70%-130%	
2,5-dibromotoluene #2	116	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	300	ug/kg
Ethylbenzene	ND	300	ug/kg
Methyl tert-butyl ether (MTBE)	ND	60	ug/kg
Naphthalene	2300	300	ug/kg
Toluene	ND	300	ug/kg
m,p-Xylenes	ND	300	ug/kg
o-Xylene	ND	300	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 16 Sample Description: 841061130-16

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:6.2Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50999 Lab Data File: J28559.D;M32629.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	20
71-43-2	Benzene	ND	4.9
108-86-1	Bromobenzene	ND	4.9
74-97-5	Bromochloromethane	ND	4.9
75-27-4	Bromodichloromethane	ND	4.9
75-25-2	Bromoform	ND	4.9
74-83-9	Bromomethane	ND	9.8
78-93-3	2-Butanone (MEK)	ND	9.8
104-51-8	n-Butylbenzene	ND	4.9
135-98-8	sec-Butylbenzene	ND	4.9
98-06-6	tert-Butylbenzene	ND	4.9
75-15-0	Carbon disulfide	ND	4.9
56-23-5	Carbon tetrachloride	ND	4.9
108-90-7	Chlorobenzene	ND	4.9
75-00-3	Chloroethane	ND	9.8
67-66-3	Chloroform	ND	4.9
74-87-3	Chloromethane	ND	9.8
95-49-8	2-Chlorotoluene	ND	4.9
106-43-4	4-Chlorotoluene	ND	4.9
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	4.9
124-48-1	Dibromochloromethane	ND	4.9
106-93-4	1,2-Dibromoethane (EDB)	ND	2.9
74-95-3	Dibromomethane	ND	4.9
95-50-1	1,2-Dichlorobenzene	ND	4.9
541-73-1	1,3-Dichlorobenzene	ND	4.9
106-46-7	1,4-Dichlorobenzene	ND	4.9
75-71-8	Dichlorodifluoromethane	ND	9.8
75-34-3	1,1-Dichloroethane	ND	4.9
107-06-2	1,2-Dichloroethane	ND	4.9
75-35-4	1,1-Dichloroethene	ND	4.9
156-59-2	cis-1,2-Dichloroethene	ND	4.9
156-60-5	trans-1,2-Dichloroethene	ND	4.9
78-87-5	1,2-Dichloropropane	ND	4.9
142-28-9	1,3-Dichloropropane	ND	4.9
590-20-7	2,2-Dichloropropane	ND	4.9
563-58-6	1,1-Dichloropropene	ND	4.9
10061-01-5	cis-1,3-Dichloropropene	ND	4.9
10061-02-6	trans-1,3-Dichloropropene	ND	4.9
60-29-7	Diethyl ether	ND	9.8
123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 16 (continued) Sample Description: 841061130-16

Date Collected:11/30/2006Matrix:SolidDate Received:12/1/2006Percent Moisture:6.2Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By:Dilution Factor:1Method:8260BSoil Extract Volume:

QC Batch#: 50999 Lab Data File: J28559.D;M32629.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	50
100-41-4	Ethylbenzene	ND	4.9
87-68-3	Hexachlorobutadiene	ND	4.9
591-78-6	2-Hexanone	ND	9.8
98-82-8	Isopropylbenzene	ND	4.9
99-87-6	4-Isopropyltoluene	ND	4.9
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	4.9
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	9.8
75-09-2	Methylene chloride	ND	4.9
91-20-3	Naphthalene	ND	4.9
103-65-1	n-Propylbenzene	ND	4.9
100-42-5	Styrene	ND	4.9
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	50
109-99-9	Tetrahydrofuran	ND	4.9
96-18-4	1,2,3-Trichloropropane	ND	4.9
630-20-6	1,1,1,2-Tetrachloroethane	ND	4.9
79-34-5	1,1,2,2-Tetrachloroethane	ND	4.9
127-18-4	Tetrachloroethene (PCE)	18	4.9
108-88-3	Toluene	ND	4.9
87-61-6	1,2,3-Trichlorobenzene	ND	4.9
120-82-1	1,2,4-Trichlorobenzene	ND	4.9
71-55-6	1,1,1-Trichloroethane	17	4.9
79-00-5	1,1,2-Trichloroethane	ND	4.9
79-01-6	Trichloroethene (TCE)	44	4.9
75-69-4	Trichlorofluoromethane	ND	9.8
95-63-6	1,2,4-Trimethylbenzene	ND	4.9
108-67-8	1,3,5-Trimethylbenzene	ND	4.9
75-01-4	Vinyl chloride	ND	9.8
95-47-6	o-Xylene	ND	4.9
	m,p-Xylenes	ND	4.9
Surrogate	Recovery	Limits	
Bromofluorobenzene	64%	78%-111%	
4.6.51.11.11	10.45		

Surrogate	Recovery	Limits
Bromofluorobenzene	64%	78%-111%
1,2-Dichloroethane-d4	104%	91%-114%
Toluene-d8	119%	86%-115%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	17	Sample Description:	841061130-17
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Analyzed:	12/05/06	Percent Moisture:	9.3
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6100	ug/kg
C9-C12 Aliphatics**	50	ND	6100	ug/kg
C9-C10 Aromatics***	50	ND	6100	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range	
2,5-dibromotoluene	99	70%-130%	
2,5-dibromotoluene #2	109	70%-130%	

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	310	ug/kg
Ethylbenzene	ND	310	ug/kg
Methyl tert-butyl ether (MTBE)	ND	61	ug/kg
Naphthalene	ND	310	ug/kg
Toluene	ND	310	ug/kg
m,p-Xylenes	ND	310	ug/kg
o-Xylene	ND	310	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 17 Sample Description: 841061130-17

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 9.3

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/06/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50999 Lab Data File: J28560.D;M32630.D

Units: ug/kg

For-64-1	CAS No.	Parameter	Result	DL
108.86-1 Bromobenzene	67-64-1	Acetone	ND	18
74-97-5 Bromochloromethane ND 4.6 75-27-4 Bromochloromethane ND 4.6 75-27-2 Bromoform ND 4.6 74-83-9 Bromomethane ND 9.2 78-93-3 2-Butanone (MEK) ND 9.2 104-51-8 n-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 56-23-5 Carbon tetrachloride ND 4.6 108-90-7 Chlorobenzene ND 4.6 75-00-3 Chloroform ND 4.6 75-00-3 Chloroform ND 4.6 74-87-3 Chloromethane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl cher (DIPE) ND 4.6 108-20-3 </td <td>71-43-2</td> <td>Benzene</td> <td>ND</td> <td>4.6</td>	71-43-2	Benzene	ND	4.6
75-27-4 Bromodichloromethane ND 4.6 75-25-2 Bromoform ND 4.6 74-83-9 Bromomethane ND 9.2 78-93-3 2-Butanone (MEK) ND 9.2 104-51-8 n-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 98-06-6 terr-Butylbenzene ND 4.6 98-06-6 terr-Butylbenzene ND 4.6 75-15-0 Carbon disulfide ND 4.6 56-23-5 Carbon tetrachloride ND 4.6 108-90-7 Chloroforme ND 4.6 75-00-3 Chloroform ND 4.6 74-87-3 Chloroform ND 4.6 74-87-3 Chloroforme ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 4.6 104-34 4.5 1.2-Dibromoethane ND 4.6	108-86-1	Bromobenzene	ND	4.6
75-25-2 Bromoferm ND 4.6 74-83-9 Bromomethane ND 9.2 78-93-3 2-Butanone (MEK) ND 9.2 104-51-8 n-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 75-15-0 Carbon disulfide ND 4.6 56-23-5 Carbon disulfide ND 4.6 75-00-3 Chlorotenzene ND 4.6 75-00-3 Chlorotene ND 4.6 75-00-3 Chlorotofum ND 4.6 75-00-3 Chlorotofum ND 4.6 75-48-3 Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 106-28-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 106-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6	74-97-5	Bromochloromethane	ND	4.6
74-83-9 Bromomethane ND 9.2 78-93-3 2-Butanone (MEK) ND 9.2 104-51-8 n-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 75-15-0 Carbon disulfide ND 4.6 108-90-7 Chlorobenzene ND 4.6 108-90-7 Chlorobenzene ND 4.6 75-00-3 Chloroform ND 9.2 67-66-3 Chloroform ND 4.6 74-87-3 Chlorofoluene ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 104-48-1 Dibromochloromethane ND 4.6 105-93-4 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6	75-27-4	Bromodichloromethane	ND	4.6
78-93-3 2-Butanone (MEK) ND 9.2 104-51-8 n-Butylbenzene ND 4.6 135-98-8 sec-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 98-06-6 tert-Butylbenzene ND 4.6 75-15-0 Carbon disulfide ND 4.6 108-90-7 Chlorobenzene ND 4.6 108-90-7 Chlorobenzene ND 4.6 75-00-3 Chlorotenane ND 9.2 67-66-3 Chlorotenane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 50 96-12-8 1,2-Dibromo-1ane ND 4.6 109-30-4 1,2-Dibromochloromethane ND 4.6 109-30-4 1,2-Dibromochlane EDB ND 2.8 74-95-3 Dibromochloromethane ND 4.6 105-31-1 1,3-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 75-71-8 Dichlorotenane ND 4.6 75-35-4 1,1-Dichlorotenane ND 4.6 75-35-4 1,1-Dichlorotenane ND 4.6 156-60-5 trans-1,2-Dichlorotenee ND 4.6 156-60-5 trans-1,2-Dichlorotenee ND 4.6 168-28-9 1,3-Dichlorotenee ND 4.6 169-28-7 Diethoropopane ND 4.6 10061-01-5 cis-1,3-Dichloropopane ND 4.6 10061-01-5 cis-1,3-Dichloropopane ND 4.6 10061-01-5 Diethyl chere ND 4.6 10061-01-5 Diethyl cher	75-25-2	Bromoform	ND	4.6
104-51-8	74-83-9	Bromomethane	ND	9.2
135-98-8 sec-Butylbenzene	78-93-3	2-Butanone (MEK)	ND	9.2
135-98-8 sec-Butylbenzene	104-51-8	n-Butylbenzene	ND	4.6
98-06-6 tert-Butylbenzene ND 4.6 75-15-0 Carbon disulfide ND 4.6 56-23-5 Carbon tetrachloride ND 4.6 108-90-7 Chlorobenzene ND 4.6 75-00-3 Chlorochane ND 9.2 67-66-3 Chloroform ND 4.6 74-87-3 Chlorochune ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 4.6 12-8 1,2-Dichlorochane	135-98-8	· · · · · · · · · · · · · · · · · · ·	ND	4.6
56-23-5 Carbon tetrachloride ND 4.6 108-90-7 Chlorobenzene ND 4.6 75-00-3 Chlorotethane ND 9.2 67-66-3 Chloroform ND 4.6 74-87-3 Chlorotethane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 106-93-4 1,2-Dibromo-dhoromethane ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 95-34-3 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 107-06-2 1,2-Dichlorotethane ND	98-06-6		ND	4.6
108-90-7	75-15-0	Carbon disulfide	ND	4.6
75-00-3 Chloroethane ND 9.2 67-66-3 Chloroform ND 4.6 74-87-3 Chloromethane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromochloromethane (EDB) ND 2.8 74-95-3 Dibromomethane (EDB) ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloropropane	56-23-5	Carbon tetrachloride	ND	4.6
67-66-3 Chloroform ND 4.6 74-87-3 Chloromethane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromo-s-chloroperopane (DBCP) ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,2-Dichlorothane ND 4.6 107-06-2 1,2-Dichlorothene ND 4.6 156-69-2 cis-1,2-Dichloropr	108-90-7	Chlorobenzene	ND	4.6
67-66-3 Chloroform ND 4.6 74-87-3 Chloromethane ND 9.2 95-49-8 2-Chlorotoluene ND 4.6 106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromo-s-chloroperopane (DBCP) ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,2-Dichlorothane ND 4.6 107-06-2 1,2-Dichlorothene ND 4.6 156-69-2 cis-1,2-Dichloropr		Chloroethane	ND	9.2
95:49-8 2-Chlorotoluene ND 4.6 106:43-4 4-Chlorotoluene ND 4.6 108:20-3 Di-isopropyl ether (DIPE) ND 50 96:12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroptopane ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 78-87-5 1,2-Dichloroprop			ND	4.6
106-43-4	74-87-3	Chloromethane	ND	9.2
106-43-4 4-Chlorotoluene ND 4.6 108-20-3 Di-isopropyl ether (DIPE) ND 50 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroptopane ND 4.6 142-28-9 1,3-Dichloroptropane ND 4.6 590-20-7 2,2-Dichloroptropane ND 4.6 563-58-6 1,1-	95-49-8	2-Chlorotoluene	ND	4.6
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7	106-43-4	4-Chlorotoluene		4.6
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 4.6 124-48-1 Dibromochloromethane ND 4.6 106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 541-73-1 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6	108-20-3	Di-isopropyl ether (DIPE)	ND	50
106-93-4 1,2-Dibromoethane (EDB) ND 2.8 74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	96-12-8		ND	4.6
74-95-3 Dibromomethane ND 4.6 95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	124-48-1	Dibromochloromethane	ND	4.6
95-50-1 1,2-Dichlorobenzene ND 4.6 541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	106-93-4	1,2-Dibromoethane (EDB)	ND	2.8
541-73-1 1,3-Dichlorobenzene ND 4.6 106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	74-95-3	Dibromomethane	ND	4.6
106-46-7 1,4-Dichlorobenzene ND 4.6 75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	95-50-1	1,2-Dichlorobenzene	ND	4.6
75-71-8 Dichlorodifluoromethane ND 9.2 75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	541-73-1	1,3-Dichlorobenzene	ND	4.6
75-34-3 1,1-Dichloroethane ND 4.6 107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	106-46-7	1,4-Dichlorobenzene	ND	4.6
107-06-2 1,2-Dichloroethane ND 4.6 75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	75-71-8	Dichlorodifluoromethane	ND	9.2
75-35-4 1,1-Dichloroethene ND 4.6 156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	75-34-3	1,1-Dichloroethane	ND	4.6
156-59-2 cis-1,2-Dichloroethene ND 4.6 156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	107-06-2	1,2-Dichloroethane	ND	4.6
156-60-5 trans-1,2-Dichloroethene ND 4.6 78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	75-35-4	1,1-Dichloroethene	ND	4.6
78-87-5 1,2-Dichloropropane ND 4.6 142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	156-59-2	cis-1,2-Dichloroethene	ND	4.6
142-28-9 1,3-Dichloropropane ND 4.6 590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	156-60-5	trans-1,2-Dichloroethene	ND	4.6
590-20-7 2,2-Dichloropropane ND 4.6 563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	78-87-5	1,2-Dichloropropane	ND	4.6
563-58-6 1,1-Dichloropropene ND 4.6 10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	142-28-9	1,3-Dichloropropane	ND	4.6
10061-01-5 cis-1,3-Dichloropropene ND 4.6 10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	590-20-7	2,2-Dichloropropane	ND	4.6
10061-02-6 trans-1,3-Dichloropropene ND 4.6 60-29-7 Diethyl ether ND 9.2	563-58-6	1,1-Dichloropropene	ND	4.6
60-29-7 Diethyl ether ND 9.2	10061-01-5	cis-1,3-Dichloropropene	ND	4.6
·	10061-02-6	trans-1,3-Dichloropropene	ND	4.6
123-91-1 1,4-Dioxane ND 18	60-29-7	Diethyl ether	ND	9.2
	123-91-1	1,4-Dioxane	ND	18

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 17 (continued) Sample Description: 841061130-17

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 9.3

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 12/06/06 By: GP Dilution Factor: 1

Method: 8260B Soil Extract Volume:

QC Batch#: 50999 Lab Data File: J28560.D;M32630.D

Units: ug/kg

CAS No.	Parameter	Res	sult	DL
	Ethyl tertiary-butyl ether (EtBE)		ND	50
100-41-4	Ethylbenzene		ND	4.6
87-68-3	Hexachlorobutadiene		ND	4.6
591-78-6	2-Hexanone		ND	9.2
98-82-8	Isopropylbenzene		ND	4.6
99-87-6	4-Isopropyltoluene		ND	4.6
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	4.6
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	9.2
75-09-2	Methylene chloride		ND	4.6
91-20-3	Naphthalene		ND	4.6
103-65-1	n-Propylbenzene		ND	4.6
100-42-5	Styrene		ND	4.6
994-05-8	Tertiary-amyl methyl ether (TAME)		ND	50
109-99-9	Tetrahydrofuran		ND	4.6
96-18-4	1,2,3-Trichloropropane		ND	4.6
630-20-6	1,1,1,2-Tetrachloroethane		ND	4.6
79-34-5	1,1,2,2-Tetrachloroethane		ND	4.6
127-18-4	Tetrachloroethene (PCE)		26	4.6
108-88-3	Toluene		ND	4.6
87-61-6	1,2,3-Trichlorobenzene		ND	4.6
120-82-1	1,2,4-Trichlorobenzene		ND	4.6
71-55-6	1,1,1-Trichloroethane		ND	4.6
79-00-5	1,1,2-Trichloroethane		ND	4.6
79-01-6	Trichloroethene (TCE)		24	4.6
75-69-4	Trichlorofluoromethane		ND	9.2
95-63-6	1,2,4-Trimethylbenzene		ND	4.6
108-67-8	1,3,5-Trimethylbenzene		ND	4.6
75-01-4	Vinyl chloride		ND	9.2
95-47-6	o-Xylene		ND	4.6
	m,p-Xylenes		ND	4.6
Surrogate	Recovery	Limits		
Bromofluorobenzene	90%	78%-111%	<u> </u>	
1,2-Dichloroethane-d4	108%	91%-114%		
Toluene-d8	90%	86%-115%		

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 18 Sample Description: 841061130-18

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Mois

Date Received:12/1/2006Percent Moisture:N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By: GPDilution Factor:50Method:8260BSoil Extract Volume:

QC Batch#: 51012 Lab Data File: M32508.D;M32617.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	250
71-43-2	Benzene	ND	50
108-86-1	Bromobenzene	ND	50
74-97-5	Bromochloromethane	ND	50
75-27-4	Bromodichloromethane	ND	50
75-25-2	Bromoform	ND	50
74-83-9	Bromomethane	ND	50
78-93-3	2-Butanone (MEK)	ND	250
104-51-8	n-Butylbenzene	ND	50
135-98-8	sec-Butylbenzene	ND	50
98-06-6	tert-Butylbenzene	ND	50
75-15-0	Carbon disulfide	ND	50
56-23-5	Carbon tetrachloride	ND	50
108-90-7	Chlorobenzene	ND	50
75-00-3	Chloroethane	ND	50
67-66-3	Chloroform	ND	50
74-87-3	Chloromethane	ND	50
95-49-8	2-Chlorotoluene	ND	50
106-43-4	4-Chlorotoluene	ND	50
108-20-3	Di-isopropyl ether (DIPE)	ND	50
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	50
124-48-1	Dibromochloromethane	ND	50
106-93-4	1,2-Dibromoethane (EDB)	ND	50
74-95-3	Dibromomethane	ND	50
95-50-1	1,2-Dichlorobenzene	ND	50
541-73-1	1,3-Dichlorobenzene	ND	50
106-46-7	1,4-Dichlorobenzene	ND	50
75-71-8	Dichlorodifluoromethane	ND	50
75-34-3	1,1-Dichloroethane	ND	50
107-06-2	1,2-Dichloroethane	ND	50
75-35-4	1,1-Dichloroethene	ND	50
156-59-2	cis-1,2-Dichloroethene	ND	50
156-60-5	trans-1,2-Dichloroethene	ND	50
78-87-5	1,2-Dichloropropane	ND	50
142-28-9	1,3-Dichloropropane	ND	50
590-20-7	2,2-Dichloropropane	ND	50
563-58-6	1,1-Dichloropropene	ND	50
10061-01-5	cis-1,3-Dichloropropene	ND	50
10061-02-6	trans-1,3-Dichloropropene	ND	50
60-29-7	Diethyl ether	ND	50
123-91-1	1,4-Dioxane	ND	1000

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 18 (continued) Sample Description: 841061130-18

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: N/A

Date Extracted:By:Sample Weight/Volume:Date Analyzed:12/06/06By: GPDilution Factor:50Method:8260BSoil Extract Volume:

QC Batch#: 51012 Lab Data File: M32508.D;M32617.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	50
100-41-4	Ethylbenzene	ND	50
87-68-3	Hexachlorobutadiene	ND	50
591-78-6	2-Hexanone	ND	250
98-82-8	Isopropylbenzene	ND	50
99-87-6	4-Isopropyltoluene	ND	50
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	50
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	250
75-09-2	Methylene chloride	ND	50
91-20-3	Naphthalene	ND	50
103-65-1	n-Propylbenzene	ND	50
100-42-5	Styrene	ND	50
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	50
109-99-9	Tetrahydrofuran	ND	50
96-18-4	1,2,3-Trichloropropane	ND	50
630-20-6	1,1,1,2-Tetrachloroethane	ND	50
79-34-5	1,1,2,2-Tetrachloroethane	ND	50
127-18-4	Tetrachloroethene (PCE)	ND	50
108-88-3	Toluene	ND	50
87-61-6	1,2,3-Trichlorobenzene	ND	50
120-82-1	1,2,4-Trichlorobenzene	ND	50
71-55-6	1,1,1-Trichloroethane	ND	50
79-00-5	1,1,2-Trichloroethane	ND	50
79-01-6	Trichloroethene (TCE)	ND	50
75-69-4	Trichlorofluoromethane	ND	50
95-63-6	1,2,4-Trimethylbenzene	ND	50
108-67-8	1,3,5-Trimethylbenzene	ND	50
75-01-4	Vinyl chloride ND		50
95-47-6			50
	m,p-Xylenes	ND	50
Surrogate	Recovery Lin	nits	
Promofluorobonzono	QQ0/4 Q70	0/ 1050/	

Surrogate	Recovery	Limits
Bromofluorobenzene	88%	87%-105%
1,2-Dichloroethane-d4	101%	91%-109%
Toluene-d8	105%	92%-105%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841061130-01
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	5.8
Date Analyzed:	12/09/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	40000	10000	ug/kg
C11-C22 Aromatics*	1	100000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	52	40%-140%
2-Bromonaphthalene	81	40%-140%
2-Fluorobiphenyl	84	40%-140%
o-Terphenyl	59	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	270	100	ug/kg
Anthracene	390	100	ug/kg
Benzo[a]anthracene	1200	100	ug/kg
Benzo[a]pyrene	1100	100	ug/kg
Benzo[b]fluoranthene	1700	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	600	100	ug/kg
Chrysene	1600	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	2200	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	990	100	ug/kg
Pyrene	2200	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 1 Sample Description: 841061130-01

Date Collected: 11/30/2006 Date Received: 12/1/2006

Date Extracted: 12/08/06 By: AKB
Date Analyzed: 12/11/06 By: LM

Method: 8082 QC Batch#: 51132 Units: ug/kg Matrix: Solid Percent Moisture: 5.8

Sample Weight/Volume: 30.19 g

Dilution Factor: 1 Extract Volume: 2 Lab Data File: 4120827.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	54%	30%-150%
Decachlorobiphenyl	53%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841061130-02
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	11.0
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	14000	10000	ug/kg
C11-C22 Aromatics*	1	16000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	50	40%-140%
2-Bromonaphthalene	71	40%-140%
2-Fluorobiphenyl	75	40%-140%
o-Terphenyl	58	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	ND	100	ug/kg
Anthracene	ND	100	ug/kg
Benzo[a]anthracene	ND	100	ug/kg
Benzo[a]pyrene	ND	100	ug/kg
Benzo[b]fluoranthene	ND	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	ND	100	ug/kg
Chrysene	ND	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	ND	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	ND	100	ug/kg
Pyrene	ND	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 Sample Description: 841061130-02

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 11.0

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.26 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120828.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15
Surrogate	Recover	ry Limits	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	80%	30%-150%
Decachlorobiphenyl	70%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841061130-03
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	10.7
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	17000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	59	40%-140%
2-Bromonaphthalene	68	40%-140%
2-Fluorobiphenyl	72	40%-140%
o-Terphenyl	52	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	740	110	ug/kg
Acenaphthylene	240	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	490	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	350	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 Sample Description: 841061130-03

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 10.7

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120829.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	69%	30%-150%
Decachlorobiphenyl	91%	30%-150%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841061130-04
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	15.8
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	32000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	54	40%-140%
2-Bromonaphthalene	62	40%-140%
2-Fluorobiphenyl	66	40%-140%
o-Terphenyl	46	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	340	110	ug/kg
Anthracene	320	110	ug/kg
Benzo[a]anthracene	1000	110	ug/kg
Benzo[a]pyrene	1000	110	ug/kg
Benzo[b]fluoranthene	1400	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	540	110	ug/kg
Chrysene	1200	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	2100	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	1100	110	ug/kg
Pyrene	2000	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 Sample Description: 841061130-04

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 15.8

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.04 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120830.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	16
11104-28-2	Aroclor 1221	ND	16
11141-16-5	Aroclor 1232	ND	16
53469-21-9	Aroclor 1242	ND	16
12672-29-6	Aroclor 1248	ND	16
11097-69-1	Aroclor 1254	ND	16
11096-82-5	Aroclor 1260	ND	16
Surrogate	Recovery	Limits	

SurrogateRecoveryLimitsTetrachloro-m-xylene63%30%-150%Decachlorobiphenyl68%30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	5	Sample Description:	841061130-05
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	14.7
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	110000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	63	40%-140%
2-Bromonaphthalene	87	40%-140%
2-Fluorobiphenyl	74	40%-140%
o-Terphenyl	58	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	210	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	1300	110	ug/kg
Anthracene	2000	110	ug/kg
Benzo[a]anthracene	4400	110	ug/kg
Benzo[a]pyrene	3900	110	ug/kg
Benzo[b]fluoranthene	5600	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	1900	110	ug/kg
Chrysene	4500	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	8200	110	ug/kg
Fluorene	390	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	3500	110	ug/kg
Pyrene	6200	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841061130-05

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 14.7

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.26 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#:

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120831.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	69%	30%-150%
Decachlorobiphenyl	74%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	6	Sample Description:	841061130-06
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	12.3
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	20000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	69	40%-140%
2-Bromonaphthalene	79	40%-140%
2-Fluorobiphenyl	81	40%-140%
o-Terphenyl	64	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	350	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	200	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	280	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	320	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	200	110	ug/kg
Pyrene	310	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 6 Sample Description: 841061130-06

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moist

Units: ug/kg

Date Received: 12/1/2006 Percent Moisture: 12.3

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.29 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120834.D

CAS No. Parameter Result DL Aroclor 1016 12674-11-2 ND 15 11104-28-2 Aroclor 1221 ND 15 11141-16-5 Aroclor 1232 ND 15 Aroclor 1242 15 53469-21-9 ND Aroclor 1248 12672-29-6 ND 15 11097-69-1 Aroclor 1254 ND 15 Aroclor 1260 15 11096-82-5 ND

SurrogateRecoveryLimitsTetrachloro-m-xylene75%30%-150%Decachlorobiphenyl76%30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	7	Sample Description:	841061130-07
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	12.3
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	40000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	57	40%-140%
2-Bromonaphthalene	73	40%-140%
2-Fluorobiphenyl	75	40%-140%
o-Terphenyl	48	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	230	110	ug/kg
Anthracene	200	110	ug/kg
Benzo[a]anthracene	330	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	560	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	170	110	ug/kg
Chrysene	640	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	380	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	150	110	ug/kg
Pyrene	470	110	ug/kg

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Matrix: Solid

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 Sample Description: 841061130-07

Date Collected: 11/30/2006 Date Received: 12/1/2006

Date Received: 12/1/2006 Percent Moisture: 12.3

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.39 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51132

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4120835.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15
Surrogate	Recovery	Limits	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	78%	30%-150%
Decachlorobiphenyl	77%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	8	Sample Description:	841061130-08
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	11.7
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	61000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	55	40%-140%
2-Bromonaphthalene	77	40%-140%
2-Fluorobiphenyl	79	40%-140%
o-Terphenyl	51	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	250	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	310	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	590	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	410	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	190	110	ug/kg
Pyrene	410	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 Sample Description: 841061130-08

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 11.7

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.19 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121122.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15
Surrogate	Recover	ry Limits	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	58%	30%-150%
Decachlorobiphenyl	105%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	9	Sample Description:	841061130-09
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	5.6
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	ND	10000	ug/kg
C11-C22 Aromatics*	1	33000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	62	40%-140%
2-Bromonaphthalene	68	40%-140%
2-Fluorobiphenyl	73	40%-140%
o-Terphenyl	53	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	220	100	ug/kg
Acenaphthylene	ND	100	ug/kg
Anthracene	970	100	ug/kg
Benzo[a]anthracene	1100	100	ug/kg
Benzo[a]pyrene	ND	100	ug/kg
Benzo[b]fluoranthene	ND	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	ND	100	ug/kg
Chrysene	1200	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	2300	100	ug/kg
Fluorene	230	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	2300	100	ug/kg
Pyrene	2000	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 9 Sample Description: 841061130-09

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 5.6

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.25 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51179

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121123.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14
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Surrogate	Recovery	Limits
Tetrachloro-m-xylene	79%	30%-150%
Decachlorobiphenyl	93%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	10	Sample Description:	841061130-10
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	8.1
Date Analyzed:	12/11/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	40	40%-140%
2-Bromonaphthalene	90	40%-140%
2-Fluorobiphenyl	88	40%-140%
o-Terphenyl	41	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Matrix: Solid

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 10 Sample Description: 841061130-10

Date Collected: 11/30/2006 Date Received: 12/1/2006

Units: ug/kg

Date Received: 12/1/2006 Percent Moisture: 8.1

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.62 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121124.D

CAS No. Parameter Result DL Aroclor 1016 12674-11-2 ND 14 11104-28-2 Aroclor 1221 ND 14 11141-16-5 Aroclor 1232 ND 14 Aroclor 1242 53469-21-9 ND 14 Aroclor 1248 14 12672-29-6 ND 11097-69-1 Aroclor 1254 ND 14 Aroclor 1260 14 11096-82-5 ND

SurrogateRecoveryLimitsTetrachloro-m-xylene71%30%-150%Decachlorobiphenyl78%30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	11	Sample Description:	841061130-11
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	6.6
Date Analyzed:	12/09/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	24000	11000	ug/kg
C11-C22 Aromatics*	1	92000	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	61	40%-140%
2-Bromonaphthalene	67	40%-140%
2-Fluorobiphenyl	76	40%-140%
o-Terphenyl	46	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	260	110	ug/kg
Anthracene	340	110	ug/kg
Benzo[a]anthracene	560	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	810	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	940	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	470	110	ug/kg
Pyrene	990	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II
PL Sample No: 11 Sample Description: 841061130-11

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moist

Date Received: 12/1/2006 Percent Moisture: 6.6
Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.84 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121125.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14
Surrogate	Recovery	I imite	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	64%	30%-150%
Decachlorobiphenyl	87%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	12	Sample Description:	841061130-12
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	14.7
Date Analyzed:	12/11/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	12000	ug/kg
C19-C36 Aliphatics	1	ND	12000	ug/kg
C11-C22 Aromatics*	1	ND	12000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	59	40%-140%
2-Bromonaphthalene	53	40%-140%
2-Fluorobiphenyl	55	40%-140%
o-Terphenyl	40	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	120	ug/kg
Acenaphthene	ND	120	ug/kg
Acenaphthylene	ND	120	ug/kg
Anthracene	ND	120	ug/kg
Benzo[a]anthracene	ND	120	ug/kg
Benzo[a]pyrene	ND	120	ug/kg
Benzo[b]fluoranthene	ND	120	ug/kg
Benzo[g,h,i]perylene	ND	120	ug/kg
Benzo[k]fluoranthene	ND	120	ug/kg
Chrysene	ND	120	ug/kg
Dibenz[a,h]anthracene	ND	120	ug/kg
Fluoranthene	ND	120	ug/kg
Fluorene	ND	120	ug/kg
Indeno[1,2,3-cd]pyrene	ND	120	ug/kg
Naphthalene	ND	120	ug/kg
Phenanthrene	ND	120	ug/kg
Pyrene	ND	120	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 12 Sample Description: 841061130-12

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 14.7

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.38 g

Date Analyzed: 12/11/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121126.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	15
11104-28-2	Aroclor 1221	ND	15
11141-16-5	Aroclor 1232	ND	15
53469-21-9	Aroclor 1242	ND	15
12672-29-6	Aroclor 1248	ND	15
11097-69-1	Aroclor 1254	ND	15
11096-82-5	Aroclor 1260	ND	15
Surrogate	Pacovary	Limite	_

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	65%	30%-150%
Decachlorobiphenyl	69%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	13	Sample Description:	841061130-13
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	8.6
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	ND	10000	ug/kg
C11-C22 Aromatics*	1	52000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	54	40%-140%
2-Bromonaphthalene	76	40%-140%
2-Fluorobiphenyl	71	40%-140%
o-Terphenyl	54	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	200	100	ug/kg
Acenaphthene	560	100	ug/kg
Acenaphthylene	150	100	ug/kg
Anthracene	1800	100	ug/kg
Benzo[a]anthracene	2800	100	ug/kg
Benzo[a]pyrene	2200	100	ug/kg
Benzo[b]fluoranthene	2600	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	1100	100	ug/kg
Chrysene	2800	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	6700	100	ug/kg
Fluorene	630	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	260	100	ug/kg
Phenanthrene	7600	100	ug/kg
Pyrene	6000	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 13 Sample Description: 841061130-13

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 8.6

Units: ug/kg

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.33 g

Date Analyzed: 12/12/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121129.D

CAS No. Parameter Result DL Aroclor 1016 12674-11-2 ND 14 11104-28-2 Aroclor 1221 ND 14 11141-16-5 Aroclor 1232 ND 14 Aroclor 1242 53469-21-9 ND 14 Aroclor 1248 14 12672-29-6 ND 11097-69-1 Aroclor 1254 ND 14 Aroclor 1260 14 11096-82-5 ND

SurrogateRecoveryLimitsTetrachloro-m-xylene81%30%-150%Decachlorobiphenyl85%30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	14	Sample Description:	841061130-14
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	8.9
Date Analyzed:	12/09/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	38000	10000	ug/kg
C11-C22 Aromatics*	1	81000	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	57	40%-140%
2-Bromonaphthalene	74	40%-140%
2-Fluorobiphenyl	73	40%-140%
o-Terphenyl	54	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	120	100	ug/kg
Anthracene	ND	100	ug/kg
Benzo[a]anthracene	ND	100	ug/kg
Benzo[a]pyrene	ND	100	ug/kg
Benzo[b]fluoranthene	230	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	ND	100	ug/kg
Chrysene	440	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	ND	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	ND	100	ug/kg
Pyrene	ND	100	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 14 Sample Description: 841061130-14

Date Collected: 11/30/2006 Matrix: Solid
Date Received: 12/1/2006 Percent Moisture: 8.9

Date Extracted: 12/08/06 By: AKB Sample Weight/Volume: 30.22 g

Date Analyzed: 12/12/06 By: LM

Method: 8082

QC Batch#: 51179

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121130.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	14
11104-28-2	Aroclor 1221	ND	14
11141-16-5	Aroclor 1232	ND	14
53469-21-9	Aroclor 1242	ND	14
12672-29-6	Aroclor 1248	ND	14
11097-69-1	Aroclor 1254	ND	14
11096-82-5	Aroclor 1260	ND	14

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	66%	30%-150%
Decachlorobiphenyl	87%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	15	Sample Description:	841061130-15
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	11.6
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	68	40%-140%
2-Bromonaphthalene	74	40%-140%
2-Fluorobiphenyl	80	40%-140%
o-Terphenyl	59	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 15 Sample Description: 841061130-15

Date Collected: 11/30/2006 Matrix: Solid

Date Received: 12/1/2006 Percent Moisture: 11.6

Date Extracted: 12/11/06 By: MM Sample Weight/Volume: 9.93 g

Date Analyzed: 12/12/06 By: LM

Method: 8082

QC Batch#: 51179

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121131.D

Units: ug/kg

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	46
11104-28-2	Aroclor 1221	ND	46
11141-16-5	Aroclor 1232	ND	46
53469-21-9	Aroclor 1242	ND	46
12672-29-6	Aroclor 1248	ND	46
11097-69-1	Aroclor 1254	ND	46
11096-82-5	Aroclor 1260	ND	46
Surrogate	Recovery	Limite	

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	71%	30%-150%
Decachlorobiphenyl	76%	30%-150%

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Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	16	Sample Description:	841061130-16
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	6.2
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	11000	ug/kg
C19-C36 Aliphatics	1	ND	11000	ug/kg
C11-C22 Aromatics*	1	ND	11000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	64	40%-140%
2-Bromonaphthalene	75	40%-140%
2-Fluorobiphenyl	77	40%-140%
o-Terphenyl	55	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
2-Methylnaphthalene	ND	110	ug/kg
Acenaphthene	ND	110	ug/kg
Acenaphthylene	ND	110	ug/kg
Anthracene	ND	110	ug/kg
Benzo[a]anthracene	ND	110	ug/kg
Benzo[a]pyrene	ND	110	ug/kg
Benzo[b]fluoranthene	ND	110	ug/kg
Benzo[g,h,i]perylene	ND	110	ug/kg
Benzo[k]fluoranthene	ND	110	ug/kg
Chrysene	ND	110	ug/kg
Dibenz[a,h]anthracene	ND	110	ug/kg
Fluoranthene	ND	110	ug/kg
Fluorene	ND	110	ug/kg
Indeno[1,2,3-cd]pyrene	ND	110	ug/kg
Naphthalene	ND	110	ug/kg
Phenanthrene	ND	110	ug/kg
Pyrene	ND	110	ug/kg

Page 100 of 103

Matrix: Solid

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 16 Sample Description: 841061130-16

Date Collected: 11/30/2006 Date Received: 12/1/2006

Date Received: 12/1/2006 Percent Moisture: 6.2

Date Extracted: 12/11/06 By: MM Sample Weight/Volume: 9.88 g

Date Analyzed: 12/12/06 By: LM

Method: 8082

QC Batch#: 51179

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121132.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	43
11104-28-2	Aroclor 1221	ND	43
11141-16-5	Aroclor 1232	ND	43
53469-21-9	Aroclor 1242	ND	43
12672-29-6	Aroclor 1248	ND	43
11097-69-1	Aroclor 1254	ND	43
11096-82-5	Aroclor 1260	ND	43
Surrogate	Pacovary	Limite	

SurrogateRecoveryLimitsTetrachloro-m-xylene73%30%-150%Decachlorobiphenyl80%30%-150%

Page 101 of 103

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E612052	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	17	Sample Description:	841061130-17
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/30/2006		
Date Received:	12/1/2006	Matrix:	Solid
Date Extracted:	12/04/06	Percent Moisture:	9.3
Date Analyzed:	12/08/06	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	10000	ug/kg
C19-C36 Aliphatics	1	ND	10000	ug/kg
C11-C22 Aromatics*	1	ND	10000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	56	40%-140%
2-Bromonaphthalene	81	40%-140%
2-Fluorobiphenyl	77	40%-140%
o-Terphenyl	49	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	100	ug/kg
Acenaphthene	ND	100	ug/kg
Acenaphthylene	ND	100	ug/kg
Anthracene	ND	100	ug/kg
Benzo[a]anthracene	ND	100	ug/kg
Benzo[a]pyrene	ND	100	ug/kg
Benzo[b]fluoranthene	ND	100	ug/kg
Benzo[g,h,i]perylene	ND	100	ug/kg
Benzo[k]fluoranthene	ND	100	ug/kg
Chrysene	ND	100	ug/kg
Dibenz[a,h]anthracene	ND	100	ug/kg
Fluoranthene	ND	100	ug/kg
Fluorene	ND	100	ug/kg
Indeno[1,2,3-cd]pyrene	ND	100	ug/kg
Naphthalene	ND	100	ug/kg
Phenanthrene	ND	100	ug/kg
Pyrene	ND	100	ug/kg

Page 102 of 103

Matrix: Solid

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E612052 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 17 Sample Description: 841061130-17

Date Collected: 11/30/2006 Date Received: 12/1/2006

Date Received: 12/1/2006 Percent Moisture: 9.3

Date Extracted: 12/11/06 By: MM Sample Weight/Volume: 9.85 g

Date Analyzed: 12/12/06 By: LM

Method: 8082

QC Batch#: 51179

Units: ug/kg

Dilution Factor: 1

Extract Volume: 2

Lab Data File: 4121133.D

CAS No.	Parameter	Result	DL
12674-11-2	Aroclor 1016	ND	45
11104-28-2	Aroclor 1221	ND	45
11141-16-5	Aroclor 1232	ND	45
53469-21-9	Aroclor 1242	ND	45
12672-29-6	Aroclor 1248	ND	45
11097-69-1	Aroclor 1254	ND	45
11096-82-5	Aroclor 1260	ND	45
_	_		

Surrogate	Recovery	Limits
Tetrachloro-m-xylene	56%	30%-150%
Decachlorobiphenyl	60%	30%-150%

Page 103 of 103



☐ 146 Hartford Road, Manchester, CT 06040

- □ 78 Interstate Drive, West Springfield, MA 01089

275 Promenade Street, Suite 350, Providence, RI 02908

olines to Deliver 46-2469 • www.FandO.com	☐ 56 Quarry Road, Trumbull, CT 06611☐ 1419 Richland Street, Columbia, SC 29201☐	 610 Lynndale Court, Suite E, Greenville, NC 27858 24 Madison Avenue Extension, Albany, NY 12203 	☐ 80 Washington Street, Suite 301, Poughkeepsie, NY 12601 ☐ Other

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P.O. No.: 8412	0050458-B10				///>///////////////////////////////////
Sampler's Signature:	Vtr Been	Date: 11/30/06			
Source Codes: MW=Monitoring Well SW=Surface Water X=Other	PW=Potable Water S=Soil T=Treatment Facility B=Bottom Sedime	W=Waste	Analysis Request Original de la companya de la com		
Item No. 1 2 3 4	Sample Number	Source Date Time Code Sampled Sampled			Comments
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- 56 Quarry Road, Trumbull, CT 06611
- ☐ 1419 Richland Street, Columbia, SC 29201
- □ 78 Interstate Drive, West Springfield, MA 01089
- □ 610 Lynndale Court, Suite E, Greenville, NC 27858

-see attacked Checkist

☐ 24 Madison Avenue Extension, Albany, NY 12203

275 Promenade Street, Suite 350, Providence, RI 02908

80 Washington Street, Suite 301, Poughkeepsie, NY 12601

Other ...

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SW=Surface Water	Γ=Treatment Facility Ε		W=Waste A=Air			/							Ci Ci Ci			\z\-\ _	~/ / /
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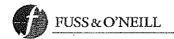
150



Quality Assurance Project Plan Phase II Site Assessment Former Nu-Style Company, Inc. Facility, Franklin, Massachusetts RFA #07011 October 2006 Revision 1.0

APPENDIX B

DATA VALIDATION COMPLETENESS CHECKLIST



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

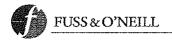
	<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREMENTS:		
Field measurement calibration records		
Groundwater field measurements (if applicable)		
Soil sampling field measurements (if applicable)		
Sediment sampling field measurements (if applicable)		
Surface water sampling field measurements (if applicable)		
Low-flow sampling field measurements (if applicable)		
Documentation of field activities		
Sample numbering and labeling		
Chain-of-Custody records		
Trip blanks		
Duplicate samples		
Equipment blanks		
Split samples (if any)		
2. LABORATORY MEASUREMENT'S:		
Trip blanks		
Instrument blanks		
Laboratory control samples		
Duplicates samples		
Equipment blanks		
Matrix spike/matrix spike duplicates		
Analysis type		
Chain-of-Custody records		
Surrogate recoveries		
Sample Project Narratives		
Split samples (if any)		
TOTAL:		
PERCENT COM	MPLETE:	%



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY FUSS & O'NEILL MODIFIED TIER II DATA VALIDATION CHECKLIST

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMITS?

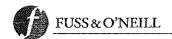
	<u>YES</u>	NO	COMMENTS
1. SAMPLING AND FIELD MEASUREMENT	S:		
Field measurement calibration records			•
pH - \pm 0.3 pH units			
S.C \pm 5% of calibration solution, within?			
calibration range			
Temperature - ± 0.5 °C			
D.O ± 5% of calibration solution			
Groundwater field measurements (if applicable)			
Water depth measured to within 0.01 ft.?			
Soil sampling field measurements (if applicable)			
OVM - ± 2 ppm			
OVA - ± 2 ppm			
Sediment sampling field measurements (if applicable)			
Descriptive information recorded?			
Surface water sampling field measurements (if applicable)			
Water depth measured to within 0.01 ft.?			
Low-flow sampling field measurements (if applicable)			
S.C ± 10%			
pH - \pm 0.2 pH units			
Temperature - ± 10%			
Turbidity - ±5 NTU			
Documentation of field activities			
Site-specific information documented in field notebook?) []		
Field data sheets completed?			
Sample numbering and labeling			
Sample numbering conforms to sample I.D. system			
identified in QAPP?			
Chain-of-Custody records			
Chain-of-Custody forms completed?	П	П	



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY FUSS & O'NEILL MODIFIED TIER II DATA VALIDATION CHECKLIST (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMITS?

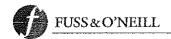
	<u>YES</u>	NO	COMMENTS
Trip blanks			
Trip blanks submitted, one per day?			
Any compounds detected in trip blanks?			
Duplicate samples			
Field duplicates performed, 1/20 samples?			
Is percent difference within 30% for all field parameters?			
Equipment blanks			
Equipment blanks submitted, one per sampling day?			
Any compounds detected in equipment blank?			
Split samples (if any)			
Split samples collected?			
Is percent difference within 30% for split samples?			
2. LABORATORY MEASUREMENTS: Trip blanks			
Trip blanks submitted, one per day?	П		
Any compounds detected in trip blanks?	П		
Instrument blanks**	$\bar{\Box}$	n	
Laboratory control samples**			
Duplicates samples**			
Equipment blanks**			
Matrix spike/matrix spike duplicates**			
Analysis type			
Chain-of-Custody records			
Surrogate recoveries**			
Sample Project Narratives			
Split samples (if any)**			
Most recent EPA WP-PE sample results**			



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

			<u>YES</u>	<u>NO</u>	<u>COMMENTS</u>
SI	DG Project Narratives				
V	olatiles Data				
a.	Sample Data				•
	• • • • • • • • • • • • • • • • • • • •				
		Ĺ.J			
	•				
					····-
	Percent solids calculations				
b.	Standards Data (all instruments)				
	Initial Calibration Data				
	RICs and Quan Reports for all Standards				
	Continuing Calibration				
	RICs and Quan Reports for all Standards				
	Internal Standard Area Summary				
c.	Raw QC Data				
	Blank Data				
	Matrix Spike Data				
	Matrix Spike Duplicate Data				
Set	mivolatiles Data				
a.	OC Summary				
	MS/MSD Summary				
	Method Blank Summary				
	Tuning and Mass Calibration				
	Tr Vo a. b.	Target Compound List (TCL) Results Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library matches Percent solids calculations b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary c. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Duplicate Data Semivolatiles Data a. QC Summary Surrogate Percent Recovery Summary MS/MSD Summary Method Blank Summary	Traffic Report Volatiles Data a. Sample Data Target Compound List (TCL) Results Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library matches Percent solids calculations b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary c. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Duplicate Data Semivolatiles Data a. QC Summary Surrogate Percent Recovery Summary Method Blank Summary	SDG Project Narratives Traffic Report Volatiles Data a. Sample Data Target Compound List (TCL) Results Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library matches Percent solids calculations b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary c. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Data Matrix Spike Duplicate Data Semivolatiles Data a. QC Summary Surrogate Percent Recovery Summary MS/MSD Summary Method Blank Summary	SDG Project Narratives Traffic Report Volatiles Data a. Sample Data Target Compound List (TCL) Results Reconstructed total ion chromatograms (RIC) for each Sample For each sample: Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library matches Percent solids calculations b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary c. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Data Matrix Spike Duplicate Data Semivolatiles Data a. QC Summary Surrogate Percent Recovery Summary MS/MSD Summary Method Blank Summary



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

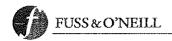
		<u>YES</u>	<u>NO</u>	COMMENTS
b.	Sample Data			
	TCL Results			
	Tentatively Identified Compounds			
	Reconstructed total ion chromatograms (RIC) for each			
	Sample			
	For each sample:			
	Raw spectra and background-subtracted mass			
	spectra of TCL compounds			
	Mass spectra of TICs with 3 best library matches			
	GPC chromatograms (if GPC performed)			
c.	Standards Data (all instruments)			
	Initial Calibration Data			
	RICs and Quan Reports for all Standards			
	Continuing Calibration			
	RICs and Quan Reports for all Standards			
	Internal Standard Areas Summary			
	Internal Standard Areas Summary			
d.	Raw QC Data			
	Decafluorotripbenylphosphine (DFTPP)			
	Blank Data			
	Matrix Spike Data			
	Matrix Spike Duplicate Data			
Mis	scellaneous Data			
	Original preparation and analysis forms or copies of preparation			
	and analysis log book pages			
	Internal sample & sample extract transfer chain-of custody			
	records			
	Screening Records			
	All instrument output, including strip charts from screening			
	activities (describe or list)			



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

I certify the abo	y that the ve analy:	e above information is a ses will be made availal	true and accurate. I further certify ole for review for seven (7) years fo	that al llowin	l labora g certif	atory results associ acation of this doc
Comple (Lab)	eted by:	(Signature)	(Printed Name/Title)			Date
	**	See laboratory Qualit	y Assurance Plan for limits.			
Comment	s: 					
Other Rec	cords (de					
Internal L	•				·· <u>·</u> ·····	
Chain-of-	Sampl	e Log-in Sheet (Lab & laneous Shipping/Re	c DC1) ceiving Records (describe or list)			



INITIAL DATE: OCTOBER 2006 REVISION DATE: OCTOBER 2006 REVISION: 1.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

	<u>YES</u>	<u>NO</u>	COMMENTS
SDG Project Narratives			
Inorganic Analysis Data Sheet			
Initial and Continuing Calibration Verification			
CRDL Standard for AA and ICP			
Blanks			
ICP Interference Check Sample			
Spike Sample Recovery			
Post Digest Spike Sample Recovery			
Duplicates			
Laboratory Control Sample			
Standard Addition Results			
ICP Serial Dilutions			
Instrument Detection Limits, Quarterly			
ICP Interelement Correction Factors, Annually			
ICP Linear Ranges Quarterly			
Preparation Log			
Analysis Run Log			
ICP Raw Data			
Furnace AA Raw Data			
Mercury Raw Data			
Percent Solids Calculations			
Digestion Logs			
EPA Shipping/Receiving Records			
(List all individual records)			
Chain-of Custody Records			
Sample Log-In sheet			
Miscellaneous Shipping/Receiving Records			
(List all individual records)		-	



INITIAL DATE: OCTOBER 2006 REVISION DATE: OCTOBER 2006 REVISION: 1.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

YES NO COMMENTS

26.	Internal Original Sample Preparation	and analysis Records	_		
	(Describe or List				
	Preparation Records				···
	Analysis Records				
	Description				
27.	Other Records (Describe or List)				
28.	Comments:				
**	See laboratory Quality Assurance Pla	n for limits.			
Comp.	eted by:				
(Lab)		(Printed Name/Title)	•	Date
associa	y that the above information is true and ted with the above analyses will be mad ation of this document.	d accurate. I further cer de available for review fo	tify the	at all labor en (7) year:	atory results s following



APPENDIX D

Phase II ESA Addendum, Former Nu-Style Property Prepared by Fuss & O'Neill February 2008

Phase II Environmental Site Assessment Addendum

Former Nu-Style Property RTN 2-0016694 87 Grove Street (Lots 22 & 27) Franklin, MA

February 2008



317 Iron Horse Way Suite 204 Providence, RI 02908



PHASE II ENVIRONMENTAL SITE ASSESSMENT ADDENDUM FORMER NU-STYLE PROPERTY 87 GROVE STREET, FRANKLIN, MA

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3.0	PHASE II ADDENDUM ACTIVITIES	8
4.0	INVESTIGATION RESULTS	11 11 11 12
5.0	EVALUATION OF ANALYTICAL RESULTS 5.1 Data Verification 5.2 Regulatory Criteria 5.3 Sediment 5.4 Soil 5.5 Groundwater	13 13 14
6.0	CONCEPTUAL SITE MODEL 6.1 Disposal Site Hydrogeology 6.2 Contaminant Sources 6.3 Nature and Extent of Contamination 6.4 Migration Pathways	15 16 17
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PHASE II ENVIRONMENTAL SITE ASSESSMENT ADDENDUM FORMER NU-STYLE PROPERTY 87 GROVE STREET, FRANKLIN, MA

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3 2 5	Parcel Number Changes	
3	Summary of Groundwater Analytical Methods	
4 5	Summary of Sediment Analytical Data and Objectives Summary of Detected Compounds in Sediment Samples	END OF REPORT 11
6 7	Summary of Soil Analytical Data and Objectives	END OF REPORT
8 9	Summary of Detected Compounds in Soil Samples Summary of Groundwater Analytical Data and Objectives Summary of Detected Compounds in Groundwater	END OF REPORT 12
9 10 11	Summary of Surface Water Analytical Data and Objectives Summary of MADEP Criteria for Detected Compounds in	END OF REPORT
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12	Groundwater Elevation Summary	END OF REPORT
<u>FIGU</u>	<u>IRES</u>	END OF REPORT
1 2	Site Location Map Site Plan	
3	Groundwater Equipotential Contour Map	
۸۵۵۲	ENDICES	END OF REPORT
WLL L	INDICES	LIND OF KLYOKI

- MADEP Bureau of Waste Site Cleanup Site Scoring Map Α
- В Soil Boring Logs and Monitoring Well Completion Reports
- С Premier Laboratory Certificates of Analysis, Fuss & O'Neill Data Verification Narratives and Certifications, and Data Validation Completeness Worksheets



1.0 INTRODUCTION

1.1 <u>Project Overview and Objectives</u>

The County of Norfolk, Massachusetts (Norfolk County) retained Fuss & O'Neill, Inc. (Fuss & O'Neill) to conduct additional Phase II Environmental Site Assessment activities (Phase II Addendum) at the former Nu-Style Company, Inc. property (the site) located at 87 Grove Street in the Town of Franklin, Massachusetts (the Town). This Phase II Addendum was conducted as part of the County Hazardous Materials and Petroleum Brownfield Assessment Programs, funded under two brownfield assessment grants from the United States Environmental Protection Agency (USEPA).

The Phase II Addendum presents the findings of the investigations performed, the conclusions drawn based on those findings, and recommendations with respect to further evaluations or other response actions that may be conducted at the site.

Previous Phase II activities conducted by Fuss & O'Neill at the site are documented in Phase II reports dated January 2007 and September 2007. Results of these investigations are also discussed herein.

The objective of the Phase II Addendum documented herein was to evaluate the nature and extent of contaminants in deep groundwater from fractured bedrock at the site and to further evaluate previously identified areas of environmental concern at the site. The Phase II Addendum scope of work was developed to determine, to the extent possible with the available resources, the absence or presence and, where applicable, the nature and extent of contaminants in environmental media, to facilitate redevelopment planning at the site, and ultimately to return the property to productive use. The Phase II Addendum was conducted in accordance with the Quality Assurance Project Plan (QAPP) Addendum Revision 3.0 dated September 2007.

1.2 <u>Assessment Planning and Approvals</u>

Prior to the commencement of initial Phase II field activities at the site, Fuss & O'Neill prepared a QAPP for review and approval by USEPA. The QAPP (Revision 0.0) was submitted to USEPA in September 2006. In October 2006, Fuss & O'Neill submitted responses to USEPA comments and questions regarding the QAPP in the form of a document titled QAPP Revision 1.0. The QAPP (Revision 1.0) was formally approved by USEPA on November 6, 2006.

QAPP Addendum (Revision 2.0) was submitted on March 14, 2007 and was formally approved by USEPA on March 20 and 23, 2007. QAPP Addendum (Revision 2.0) detailed the field and analytical scope for the collection of surface water and sediment samples as well as the closure assessment of an underground storage tank (UST).

QAPP Addendum (Revision 3.0) was submitted on September 7, 2007 and was formally approved by USEPA on October 9, 2007. QAPP Addendum (Revision 3.0) detailed the field and analytical scope for additional sediment, soil, and groundwater sampling at the site, as summarized herein.



The QAPPs, developed in accordance with the USEPA Brownfields Quality Assurance Project Plan Guidance Document, detailed the field and analytical scope and quality control procedures to be implemented during the Phase II activities. The above-referenced QAPPs are hereafter referred to collectively as the "approved QAPPs."

2.0 BACKGROUND

2.1 Site Description

The site was located at 87 Grove Street in Franklin, Massachusetts (UTM NAD83 meters: Northing 4,662,290 Easting 299,210; Lat/Long: 42°5′13.154″ N 71°25′39.790″ W). The site was identified as the Town of Franklin Tax Assessor's Map 276, Lot 22 and Lot 27. Lot 22 covered an area of approximately 9,929 square feet. Lot 27 adjoined Lot 22 to the east and was approximately 42,359 square feet in size. The site was acquired via tax title by the Town of Franklin as a result of foreclosure. Details of the site history are presented in the QAPP (Revision 1.0) and Phase I ESA, prepared by Fuss & O'Neill in 2006.

A vacant, partially dilapidated two-story building with a footprint of approximately 11,800 square feet was situated on Lot 27, and a vacant one and one-half-story building with a footprint of approximately 4,000 square feet was located on Lot 22. Mine Brook flowed westward along the southern side of the Lot 27 building and turned northward to form the western boundary of Lot 22. Mine Brook flowed generally northward to the Charles River. Unrestricted access to the subject property was provided via Grove Street and Old Grove Street.

Utilities located on the subject property include a water line located along the right-of-way known as Old Grove Street, and stormwater drainage lines located throughout the property.

No endangered species habitat, areas of critical environmental concern (ACEC) or certified vernal pools were located within 500 feet of the subject site, as depicted on the Massachusetts Department of Environmental Protection (MADEP) Bureau of Waste Site Cleanup Site Scoring Map attached as <u>Appendix A</u>.

A portion of a United States Geological Survey (USGS) topographic map depicting the location of the site is provided as <u>Figure 1</u>. A site plan, depicting the boundary of the disposal site, is provided as <u>Figure 2</u>.

2.2 Groundwater Classification

According to the Massachusetts Contingency Plan (MCP) (310 CMR 40.0932), groundwater at the subject site is classified as GW-2/GW-3. All groundwater in the Commonwealth of Massachusetts is considered a potential source of discharge to surface water and, therefore, is categorized, at a minimum, as class GW-3.

GW-2 also applies to the site because groundwater at the site is typically present at depths of less than 15 feet below grade and, at the northern portion of the site, is within 30 feet of a potentially occupied building. In addition, it is likely that regularly occupied structures will



be present at the site subsequent to redevelopment. Category GW-2 groundwater is considered a potential source of vapors of oil and/or hazardous material (OHM) to indoor air.

The site is not located within a MADEP Zone II (aquifer protection area), potentially productive aquifer, or other GW-1 inclusionary criteria (MassGIS, 2007); therefore, a classification of GW-1 does not apply to the property. The MADEP Bureau of Waste Site Cleanup Site Scoring Map is attached as <u>Appendix A</u>.

2.3 <u>Location of Public Water Sources</u>

No public water supply wells or systems are located within a one-half-mile radius of the subject site; however, a public water supply system associated with Franklin Water Department Well #2 is located just over one-half-mile to the southeast of the subject site, at Beaver Pond (Fuss & O'Neill, 2007a). This area is classified as Zone II (aquifer protection area). Based on the inferred groundwater flow direction, it is unlikely that releases that may have occurred at the subject site would have an adverse impact on groundwater quality within the aquifer protection area.

Several United States Geological Survey (USGS) wells were also located near Beaver Pond as well as within a one-half-mile radius of the subject site. USGS wells within a one-half-mile radius of the subject site are listed in <u>Table 1</u> below.

Table 1 Summary of USGS Wells

Well	Distance/Direction from Subject Site
USGS 3319020	~ 0.1 mile/East
USGS 3319051	~ 0.15 mile/North
USGS 3319013	~ 0.15 mile/East-southeast
USGS 3319068	~ 0.45 mile/Northwest
USGS 3319084	~ 0.5 mile/North-northwest

2.4 <u>Topography and Geology</u>

The topography of the site was generally flat, except at the banks of Mine Brook, where the topography dropped steeply to the river bed (USGS, 1987). The regional topography was hilly and generally drained to Mine Brook.

Surficial material at the site was mapped as loamy udorthents, which generally consist of moderately coarse-grained, deep and moderately deep, fairly well-drained soils (USDA, 2006). Fill described as sand, gravel, silt, and, in some cases, wood and brick was observed to depths of up to 14 feet below grade during drilling conducted on the site as part of the Phase II documented herein.



Bedrock beneath the site was mapped as grayish-pink to greenish-gray, equigranular to slightly porphyritic, Dedham Granite (Zen, 1983). Bedrock was encountered at the site during drilling at depths of between four and 12.5 feet below grade.

2.5 <u>Previous Environmental Investigations</u>

The following reports document the results of environmental investigations conducted at the subject site thus far:

- IES, Inc., 1990. Chapter 21-E Site Evaluation, 87 Grove Street, Franklin, MA. January 17, 1990.
- IES, Inc., 1991. Test Borings and Analysis, 87 Grove Street, Franklin, MA. July 24, 1991.
- Fuss & O'Neill, 2006. Phase I Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, May 2006.
- Fuss & O'Neill, 2007. Phase II Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, January 2007.
- Fuss & O'Neill, 2007. Phase I Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, February 2007.
- Fuss & O'Neill, 2007. UST Closure Assessment Report, Former Nu-Style Company, Inc., Franklin, Massachusetts, July 2007.
- Fuss & O'Neill, 2007. Phase II Environmental Site Assessment, Former Nu-Style Company, Inc. Facility, 87 Grove Street (Lots 22 & 27), Franklin, Massachusetts, September 2007.

2.5.1 IES, Inc. Site Investigation Reports

Portions of two reports prepared by IES, Inc. (IES) summarizing environmental investigations previously conducted on the subject property and on the parcel adjacent to the south were reviewed. The results of the investigations documented in these reports are discussed below.

January 1990

In January 1990, IES completed a report of a Chapter 21E Site Evaluation of 87 Grove Street for Home National Bank of Milford, Massachusetts. Portions of the report were available for review at the Franklin Health Department.



The IES investigation included the drilling of soil borings and the collection and analyses of soil and groundwater samples on the subject property and on the parcel adjacent to the south of Lot 27 (Lot 26). Note that the map and parcel numbers have changed since the IES investigation, as summarized in the table below.

Table 2
Parcel Number Changes

Prev	vious	Cui	rrent	Comments				
Map	Lot	Map	Lot	Comments				
72	5	276	22	Site				
72	6	276	27	Site				
72	7	276	26	Adjacent south				

IES collected soil and/or groundwater samples from five borings (B-1 through B-5) drilled on the three parcels. A figure provided by IES depicted the approximate boring locations; however because the figure was schematic and was not to scale, the precise boring locations could not be determined.

Two of the borings (B-1 and B-2) were drilled adjacent to underground storage tanks (USTs) located on Lot 26. Borings B-3 and B-5 were situated on the north side of the Lot 27 building, and boring B-4 was advanced in the exterior "barrel area" north of the Lot 22 garage. Field screening indicated the presence of trace concentrations of volatile organic compounds (VOCs) in the soil at borings B-4 and B-5; therefore, soil from the two borings from a depth of approximately five feet below grade was composited into one sample, which was analyzed for VOCs. No VOCs were detected at levels above the laboratory reporting limit. Groundwater was not encountered at these two boring locations.

Groundwater samples collected from borings B-1 and B-2 were analyzed for VOCs and were also not detected at concentrations above the laboratory reporting limit. No information regarding sample analysis for soil or groundwater collected from boring B-3 was reported; therefore, we infer that no samples were analyzed because field screening did not indicate the presence of VOCs.

IES concluded that no releases of hazardous materials or petroleum products had occurred at the subject property; however, it is Fuss & O'Neill's opinion that the IES investigation was not adequate to definitively rule out releases on the subject property.

July 1991

In July 1991, IES collected soil samples from four additional borings (B-1A through B-4A) to assess whether releases associated with USTs had occurred. As with the 1990 investigation, only portions of the July 1991 report were available for review at the Franklin Health Department. A copy of the report was also available at the Franklin Fire Department, but copies could not be made. A figure depicting the boring locations was not included with the report.

Soil encountered at the site generally consisted of fill containing loam, sand, gravel, and, in some cases, brick and cinders. Fill materials were observed to depths of up to 8.5 feet



below grade (fbg). Deeper soil consisted of very dense, fine-grained sand, silt, and gravel. Groundwater was encountered at depths of approximately 8.5 to 9 fbg. Monitoring wells were installed within the borings to allow for the collection of groundwater samples.

IES identified releases of chlorinated solvents to soil and groundwater at boring location B-4A, which was advanced downgradient of USTs at the site and north of Mine Brook. Based on the apparent vertical distribution of VOCs in soil, IES inferred that the presence of VOCs was the result of a surface release.

2.5.2 Phase I ESA Report, May 2006

A Phase I ESA, prepared by Fuss & O'Neill in May 2006, identified the following recognized environmental conditions (RECs) at the site:

- The site had a long history (at least 90 years) of manufacturing, including textiles and jewelry. Materials used and stored at the site associated with jewelry manufacturing included cyanides, metals, chlorinated solvents, and petroleum products. Additional substances associated with textile manufacturing were also likely used. Files indicated that numerous drums of hazardous waste and petroleum products were located outside of the site buildings.
- At least one UST was present on the western side of the Lot 27 building. In addition, a heating oil tank reportedly existed in an underground bunker on the same side of the building.
- A small tunnel containing slow-flowing water was present beneath the Lot 22 building. A review of mapping on file at the Town Building Department suggested that the tunnel runs, or ran in the past, from Mine Brook and beneath the Lot 27 building to the Lot 22 building. There is the potential that the tunnel was used by the former woolen mill for direct waste disposal to Mine Brook prior to the realignment of the brook in the 1960s.
- Releases of chlorinated solvents to soil and groundwater were identified on Lot 26, which abutted the site to the south. Due to the proximity of this property to the site, there is the potential for releases that occurred on this property to adversely affect groundwater quality at the site. Note that this property was owned and occupied by the same entities that owned and operated the facilities at the site; therefore, there is the potential that similar releases have occurred at the site.
- The southern portion of the site contained a pond that was filled circa 1960. The fill appeared to have been placed by a municipality. The nature and origin of the fill were not known.



2.5.3 Phase I ESA Report, January 2007

A Phase I ESA report was prepared by Fuss & O'Neill in accordance with ASTM E1527-05 dated January 2007. The January Phase I ESA report identified the same RECs that were noted in May 2006, summarized above.

2.5.4 UST Closure Assessment Report, July 2007

Fuss & O'Neill conducted oversight of the removal of a 5,000-gallon #2 heating oil UST at the site on May 1 and 2, 2007, in accordance with QAPP Addendum (Revision 2.0). Closure activities were completed by TMC Services, Inc. (TMC) of Bellingham, Massachusetts.

Following tank removal, Fuss & O'Neill collected six confirmation soil samples from the limits of the excavation. The six confirmation soil samples were submitted to a fixed-base laboratory for analysis of the Priority Pollutant 13 metals (PP13 metals), VOCs, and petroleum hydrocarbons. In accordance with 310 CMR 40.0361, the reportable concentrations in soil category 1 (RCS-1) reporting category and the MCP Method 1 Standard Application for S-1, S-2, and S-3 for GW-1, GW-2, and GW-3 areas were applied to the confirmatory soil samples obtained on the subject site. Laboratory analytical results of soil samples collected from the limits of the tank grave did not document the presence compounds at levels in excess of the applicable criteria.

2.5.5 Phase II ESA, November 2006 through May 2007

Fuss & O'Neill conducted Phase II ESA field activities at the site during November and December 2006, and again in April and May 2007. Field activities included soil, sediment, and groundwater sampling throughout the site to characterize soil and groundwater associated with the environmental concerns identified in the Phase I ESA conducted by Fuss & O'Neill in May 2006. The results of the Phase II field investigations were documented in the Phase II ESA reports prepared by Fuss & O'Neill in January and September 2007, and are summarized below:

Soil

A comparison of the soil analytical results to the RCS-1 standards indicated that pursuant to the MCP (310 CMR 40.0361), a reportable condition existed with regard to soil at the subject site.

Laboratory analytical results of soil samples collected from soil borings advanced on-site documented the presence of the following target analytes at concentrations in excess of one or more criteria:

- Metals: beryllium, lead and nickel
- VOCs: Tetrachloroethylene (PCE) and trichloroethylene (TCE)
- Polynuclear Aromatic Hydrocarbons (PAHs): benzo(a)pyrene and fluorene

Sediment

A comparison of the sediment analytical results to the MADEP Revised Sediment Screening Values, which incorporated the Threshold Effects Concentrations (TECs), and the RCS-1 reporting category, indicated that acenaphthylene, anthracene, fluoranthene, phenanthrene,



and pyrene were detected at concentrations in excess of the sediment screening values. Samples did not exceed the RCS-1 criteria.

<u>Groundwater</u>

A comparison of the analytical results in shallow overburden groundwater samples to the reportable concentrations in groundwater category 2 (RCGW-2) indicated that pursuant to the MCP, 310 CMR 40.0362, a reportable condition existed with regard to groundwater at the subject site.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of the following target analytes in shallow overburden groundwater at concentrations in excess of one or more criteria:

Metals: lead

VOCs: PCE and TCE

Surface Water

A comparison of the surface water analytical results to the MCP Method 1 Groundwater Standard Application for GW-3 areas and the USEPA Chronic Criteria Continuous Concentrations indicated that no samples contained concentrations of analytes that exceeded the reference criteria.

Based on the initial Phase II ESA results, a release notification form was submitted to the MADEP on May 10, 2007. MADEP assigned release tracking number (RTN) 2-0016694 to the site.

3.0 PHASE II ADDENDUM ACTIVITIES

Based on the results of previous Phase II activities, a Phase II Addendum scope of work was developed to assess the nature and extent of contaminants in deep groundwater from fractured bedrock at the site, and to further evaluate previously identified areas of environmental concern at the site. The Phase II Addendum scope of work for soil, groundwater, and sediment sampling was implemented in accordance with the approved OAPPs.

3.1 <u>Field Investigation Activities</u>

3.1.1 Sediment Sampling

In accordance with the approved QAPPs, a sediment sampling program was implemented for the site on October 25, 2007. Sediment samples were collected from three locations, designated as SD-5 through SD-7, from the banks of Mine Brook. The purpose for the sediment sampling was to delineate the extent of sediment containing PAHs in the vicinity of sediment sample SD-01. Refer to Figure 2 for a map of sediment sample locations.

Four sediment samples, including one field duplicate collected from location SD-5, were submitted to Premier Laboratory, LLC (Premier) of Dayville, Connecticut for analysis of semi-volatile organic compounds (SVOCs) (EPA Method 8270C). An aqueous trip blank



was not submitted for analysis because VOCs were not an established compound of concern. Dedicated sampling materials were utilized to collect the sediment samples. Therefore, an equipment blank was not collected for laboratory analysis.

3.1.2 Soil Sampling

In accordance with the approved QAPPs, a soil sampling program was conducted at the site on October 31 and November 1, 2007. Five soil borings were advanced at the site. Three shallow borings, designated B-15, MW-16, and MW-17, were advanced utilizing a hollow-stem auger (HSA) drill rig to a depth of up to 14 fbg. Two deep borings, designated MW-13 and MW-14, were advanced utilizing HSA (overburden) and air-rotary (bedrock) drilling methods. The purpose for the soil boring and soil sampling program was to further delineate the extent of soil containing VOCs and metals as well as to facilitate the installation of additional monitoring wells at the site. Refer to Figure 2 for a map of soil boring locations.

One soil sample was collected from boring MW-13 from vadose zone soil directly above the water table, and one shallow soil sample was collected from boring B-15. Three soil samples were collected from boring MW-17; one shallow soil sample and two soil samples, including one field duplicate, from vadose zone soil directly above the water table.

Five soil samples were submitted to Premier for analysis. Soil samples collected from borings MW-13 and B-15, located north of Mine Brook, were analyzed for VOCs by USEPA Method 8260B. Soil samples collected from boring MW-17, located south of Mine Brook, were analyzed for VOCs by USEPA Method 8260B, PP13 metals plus barium by USEPA Methods 6010B and 7471, and petroleum hydrocarbons by MADEP EPH and VPH methods.

Soil boring logs, depicting sample recovery amounts, material descriptions, graphic logs, soil codes, and photoionization detector (PID) field screening results are attached in Appendix B.

3.1.3 Monitoring Well Installation and Development

In accordance with the approved QAPPs, groundwater monitoring wells were installed in four of the soil borings advanced at the site. Wells were installed into and screened in bedrock at borings MW-13 and MW-14, and installed and screened into overburden at borings MW-16 and MW-17. The locations of the monitoring wells are depicted on <u>Figure 2</u>. Detailed monitoring well completion reports are included in <u>Appendix B</u>.

In accordance with the approved QAPPs, Fuss & O'Neill surveyed the relative elevations of the newly installed monitoring wells at the site on November 7, 2007. The survey was conducted relative to an assumed arbitrary vertical datum to evaluate the relative elevation and hydraulic gradient of shallow groundwater beneath the site.

In accordance with the approved QAPPs, Fuss & O'Neill developed the newly installed monitoring wells at the site on November 5, 2007. Development procedures included the repeated purging and surging of groundwater in the wells to remove fine particles and to



improve hydraulic communication between the sand filter pack and surrounding soil formation.

3.1.4 Low Flow Groundwater Sampling

On November 6 and November 7, 2007, Fuss & O'Neill collected ten groundwater samples, including two duplicate samples, from monitoring wells at the site. Groundwater samples were collected from monitoring wells MW-1, MW-13, MW-14, MW-16 and MW-17 utilizing low-flow sampling techniques, in accordance with the approved QAPPs.

Low water volume and slow recharge at wells MW-2, MW-3 and MW-5 prevented the utilization of low-flow field parameter monitoring techniques. Instead, groundwater samples were collected at low flow rates following limited purging. Due to the relatively high turbidity observed in purge water from these wells, collected groundwater samples were field filtered through a dedicated 0.45 micron cartridge filter, in accordance with Fuss & O'Neill's SOPs.

Previously installed monitoring well MW-4 could not be located visually or by metal detector during the groundwater sampling event, and may have been damaged following the previous groundwater sampling event.

Groundwater samples were submitted to Premier for analysis. The following table summarizes the analyses performed:

Table 3
Summary of Groundwater Analytical Methods

Sample ID	Well ID	Analytical Methods
06-01; 06- 02*	MW-17	VOCs (USEPA Method 8260B), Total RCRA 8 Metals (USEPA Methods
06-03	MW-1	6010B/7471), MADEP VPH/EPH
06-04	MW-14	
06-05; 06- 06*	MW-13	VOCs (USEPA Method 8260B), Total RCRA 8 Metals (USEPA Methods 6010B/7471)
07-08	MW-16	
07-09	MW-5	VOCa (LICEDA Mathad 00/OD Dissahud DCDA 0 Matala (LICEDA Mathad
07-10	MW-2	VOCs (USEPA Method 8260B, Dissolved RCRA 8 Metals (USEPA Method 6010B/7471)
07-11	MW-3	00100771119

NOTES:

Only the last four digits of the sample number are shown.

* indicates duplicate sample

Two groundwater samples collected during this assessment were duplicate samples submitted to the laboratory for quality control purposes. Duplicates were collected from the overburden aquifer (MW-17) and bedrock aquifer (MW-13). One trip blank per day of sampling was also collected and submitted to Premier for analysis of VOCs.



4.0 INVESTIGATION RESULTS

4.1 <u>Sediment Laboratory Analytical Results</u>

A summary of sediment analytical data is included in <u>Table 4</u>. The complete Premier analytical data packages and associated data verification narratives and certifications for each laboratory report are attached in <u>Appendix C</u>.

Laboratory analytical results of sediment samples collected from Mine Brook as part of this Phase II Addendum documented the presence of the following analytes in sediment at concentrations above laboratory reporting limits:

Table 5
Summary of Detected Compounds in Sediment Samples
Collected As Part of Phase II Addendum

SVOCs (USEPA Method 8270)
Acenaphthylene
Anthracene
Benzo(a)anthracene
Benzo(a)pyrene
Benzo(b)fluoranthene
Benzo(k)fluoranthene
Chrysene
Dibenzo(a,h)anthracene
Fluoranthene
Fluorene
Indeno(1,2,3-cd)pyrene
Phenanthrene
Pyrene

4.2 Surficial and Subsurface Soil Characterization

In general, the soil within soil borings advanced at the site was observed to consist of mainly fine to medium sand, with varying proportions of gravel and silt. Apparent fill material containing metal slag and coal and/or coal ash was observed in soil borings advanced north of Mine Brook, and was concentrated in the upper two feet of soil. Soil boring MW-17 ended in a soil horizon consisting predominately of silt and clay from 12 to 14 feet below grade. The silt and clay horizon may represent the sediment/water interface of the former pond that existed at this location prior to the historical application of fill material.

Bedrock encountered at the site was described as granite.

4.3 <u>Soil PID Field Screening Results</u>

During soil boring advancement, soil samples were collected throughout the soil column at each boring location for field screening for the presence of total VOCs with a PID. Field screening results indicated that total VOCs were detected in soil samples at concentrations up to 9.5 parts per million (ppm). Field screening results are included in soil boring logs attached as <u>Appendix B</u>.



4.4 Soil Laboratory Analytical Results

A summary of soil analytical data is included in <u>Table 6</u>. The complete Premier analytical data packages and associated data verification narratives and certifications for each laboratory report are attached in <u>Appendix C</u>.

Laboratory analytical results of soil samples collected from on-site soil borings documented the presence of the following analytes in soil at concentrations above laboratory reporting limits:

Table 7
Summary of Detected Compounds in Soil Samples
Collected As Part of Phase II Addendum

Metals (Method 6010)	VOCs (Method 8260)	PAHs (MADEP EPH Method)
Antimony	Acetone	Acenaphthene
Arsenic	Methyl Ethyl Ketone	Acenaphthylene
Barium	Naphthalene	Anthracene
Beryllium	Tetrachloroethene (PCE)	Benzo(a)anthracene
Cadmium	Trichloroethene (TCE)	Benzo(a)pyrene
Chromium		Benzo(b)fluoranthene
Copper		Benzo(k)fluoranthene
Lead		Chrysene
Mercury (Method 7471)		Fluoranthene
Nickel		Fluorene
Zinc		Indeno(1,2,3-cd)pyrene
		Phenanthrene
		Pyrene

4.5 <u>Groundwater Laboratory Analytical Results</u>

A summary of groundwater analytical data is included in <u>Table 8</u>. The complete Premier analytical data packages and associated data verification narratives and certifications for each laboratory report are attached in <u>Appendix C</u>.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of the following analytes in groundwater at concentrations above laboratory reporting limits:

Table 9
Summary of Detected Compounds in Groundwater Samples
Collected As Part of Phase II Addendum

Metals (Method 6010)	VOCs (Method 8260)
Barium	cis-1,2-Dichloroethene
Chromium	Methyl tert butyl ether (MTBE)
Lead	PCE
	TCE



5.0 EVALUATION OF ANALYTICAL RESULTS

5.1 Data Verification

Procedures and methodologies for the collection and analyses of soil and groundwater samples were performed consistent with the approved QAPPs and the MCP (310 CMR 40.0017). Analytical data were developed and reviewed in accordance with MADEP's Compendium of Quality Assurance and Quality Control Requirements and Performance Standards for Selected Analytical Methods (the CAM).

Fuss & O'Neill conducted modified Tier II data verification of the field and analytical data resulting from the assessment documented herein. Modified Tier II verification narratives and certifications, signed by the Fuss & O'Neill Quality Assurance/Quality Control Officer, as well as modified Tier II completeness and verification checklists are attached to each Premier report in Appendix C.

Presumptive Certainty was obtained for each data set collected as part of the Phase II Addendum investigation. Documentation was provided by Premier along with narrative summaries (Appendix C).

In general, reporting limits were low enough to allow for direct comparison to the applicable criteria. Due to matrix interferences, the reporting limit for dibenzo(a,h)anthracene in sediment sample SD-06 was unable to meet the applicable criteria. The lowest achievable reporting limit for hexachlorobutadiene in soil and groundwater samples did not meet applicable standards.

The relative percent differences calculated for the primary and duplicate sediment for several compounds in the sediment samples, several metals in the soil samples, and lead in the groundwater samples were greater than the limits established in the approved QAPPs. These differences are attributed to sample heterogeneities.

The usability of the data is not anticipated to be affected by these issues.

5.2 Regulatory Criteria

Analytical results were compared to the following regulatory criteria:

- Sediment —The MCP has not established regulatory criteria for sediment. However, the following reference criteria were utilized as screening values.
 - Massachusetts Revised Sediment Screening Values, which incorporated the TECs.
 - o MCP Method 1 S-1, S-2, and S-3 Standard Application for GW-2 and GW-3 areas.
- Soil
 - MCP Method 1 S-1, S-2, and S-3 Standard Application for GW-2 and GW-3 areas.



Groundwater

- MCP Method 1 Groundwater Standard Application for GW-2 and GW-3 areas (GW-2 and GW-3).
- Surface Water —although additional surface water samples were not collected as part
 of this Phase II Addendum, the results of previous surface water samples were
 tabulated herein and included as <u>Table 10</u>. Further discussion of surface water data
 is included in Section 6.0.
 - o MCP Method 1 Groundwater Standard Application for GW-3 areas.
 - USEPA Chronic Criteria Continuous Concentrations

Applicable soil and sediment criteria, as discussed above were tabulated for detected compounds in soil and sediment samples collected from the site. These standards are summarized in Table 11.

5.3 Sediment

Sediment analytical results were compared to the sediment screening values and applicable soil criteria. A summary of sediment analytical results for all samples collected by Fuss & O'Neill is included as <u>Table 4</u>.

Laboratory analytical results of sediment samples collected from Mine Brook documented the presence of several EPH or SVOC compounds in samples SD-01, SD-05, SD-06, and SD-07 at levels in excess of the sediment screening values. The detected compounds are collectively identified as PAHs. These sediment samples were collected in close proximity to each other and were collected from the western portion of the site in the vicinity of sediment sample SD-01.

5.4 Soil

In accordance with 310 CMR 40.0361, the MCP Method 1 S-1, S-2, and S-3 Standard Application for GW-2 and GW-3 areas were applied to soil samples obtained on the subject site. A summary of soil analytical results for all samples collected by Fuss & O'Neill is included as <u>Table 6</u>.

Concentrations of metals, EPH, and VPH compounds in the soil samples collected from soil borings advanced as part of the Phase II Addendum did not exceed the applicable criteria. The concentration of PCE in the soil sample collected from boring B-15 at two to four fbg exceeded the S-1/GW-2, S-1/GW-3, S-2/GW-2, and S-3/GW-2 standards, and the concentration of TCE in the sample exceeded the S-1/GW-2, S-2/GW-2, and S-3/GW-2 standards.

Overall, the concentration of PCE and TCE exceeded the applicable standards in soil samples collected from borings B-04, B-06, and B-15. Additionally, the concentration of TCE in the soil sample collected from boring B-10 exceeded applicable standards. These soil samples were collected from the vadose zone to as deep as six fbg. These results suggest that these compounds were introduced to the subsurface via surficial releases.

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The concentration of lead in soil samples collected from boring MW-05 up to seven fbg exceeded the applicable criteria. The concentration of beryllium in the soil sample collected from boring B-10 exceeded the S-2/GW-2 and S-2/GW-3 standards.

5.5 <u>Groundwater</u>

In accordance with 310 CMR 40.0362, the MCP Method 1 Groundwater Standard Application for GW-2 and GW-3 areas reporting categories were applied to all groundwater samples obtained on the subject site. A summary of groundwater analytical results for all samples collected by Fuss & O'Neill is included as <u>Table 8</u>.

Laboratory analytical results of groundwater samples collected from on-site monitoring wells documented the presence of lead at levels in excess of the GW-3 criteria in samples collected from monitoring wells MW-1, MW-3, and MW-5. The VOC compounds PCE and TCE were reported at levels in excess of the GW-2 criteria in samples collected from monitoring wells MW-3, MW-4, and MW-13. Additionally, the concentration of TCE in the sample collected from monitoring well MW-16 exceeded the GW-2 standard.

6.0 CONCEPTUAL SITE MODEL

A conceptual site model (CSM) has been developed for the site based on the nature and source of the release, geologic and hydrogeologic conditions, historical site uses, current uses and foreseeable site uses. Available site data, including data presented herein and generated in previous investigations by Fuss & O'Neill, was evaluated in developing the CSM. The CSM was used to develop conclusions regarding the extent of contamination, media affected by the releases, and sufficiency of investigations. A discussion of the source, site hydrogeology, migration pathways, and the nature and extent of contamination follows.

6.1 <u>Disposal Site Hydrogeology</u>

The disposal site was underlain by fine to medium sand, with varying proportions of gravel and silt. Apparent fill material containing metal slag and coal and/or coal ash was observed in soil borings advanced north of Mine Brook, and was concentrated in the upper two feet of soil. The soil deposits extended to at least 14 fbg based on the deepest soil boring (MW-17). No wide-ranging impermeable confining layer was identified in the available soil data at the site. Bedrock composed of granite was encountered as shallow as four fbg at soil boring B-15.

Two groundwater sampling events have been conducted at the disposal site and provide data related to seasonal groundwater occurrence and flow. A summary of groundwater elevation data for monitoring events conducted during November 2007 is presented in <u>Table 12</u>. The depth to groundwater ranged from approximately 4.5 fbg to approximately 10.5 fbg across the site. Shallow groundwater flow direction was generally to the south-southwest toward Mine Brook. An equipotential contour map of shallow groundwater is presented as Figure 3.



Two monitoring wells, designated MW-13, and MW-14, were advanced and screened in bedrock at the first water-bearing fracture encountered during drilling. Based on the groundwater elevations in these wells, groundwater flow direction in the bedrock aquifer was generally to the southward in the direction of Mine Brook. Groundwater elevations at these bedrock wells compared to the elevation of shallow groundwater in overburden monitoring wells indicated an upward vertical hydraulic gradient from the bedrock aquifer to the overburden aquifer. However, the hydraulic connection of the bedrock and overburden aquifers was not established during this investigation.

Based on a review of MCP GW-1 inclusionary criteria and the Site Scoring Map, attached as <u>Appendix A</u>, groundwater at the site and the surrounding area was not used for drinking water and was not considered a potential drinking water source.

6.2 Contaminant Sources

Based on the investigations and reports summarized herein, the following releases were identified:

- Sediment: PAHs
 - o PAHs in sediment were identified in the vicinity of sediment samples SD-01, SD-05, SD-06, and SD-07 at the western portion of the site and at downstream locations of Mine Brook.
 - o The source of PAHs in sediment was attributed to historic urban filling practices at the site in which Mine Brook was channelized and former surface waters at the site were filled in; or to historic discharges from the former manufacturing facility.
- Soil: chlorinated VOCs (PCE and TCE), beryllium, and lead
 - The highest concentrations of VOCs were identified in soil samples collected proximal to the western and northern edges of the building at two general locations:
 - § In shallow soil in the vicinity of B-06, B-10, and B-15. The source of VOCs in this area may be attributed to historic surficial spills that migrated to the subsurface.
 - § In shallow soil and soil at the approximate depth of the water table at soil boring B-04. The source of VOCs in this area may be attributed to both historic surficial spills as well as a potential additional source of VOCs in groundwater at depth.
 - o The source of beryllium in the soil sample collected from boring B-10 and lead in the soil sample collected from boring MW-05 is likely attributed to natural sources and/or historic urban filling practices at the site.
- Shallow Overburden Groundwater: chlorinated VOCs (PCE and TCE) and lead



- OVOCs in shallow groundwater were identified in samples collected from monitoring wells MW-03, MW-04, and MW-16. These wells were located in the vicinity of soil containing VOCs, likely as a result of surficial spills that migrated to groundwater. The dissolved-phase VOCs in these areas indicated that VOCs migrated to groundwater as a result of the historic spills.
- Lead in shallow groundwater was identified in samples collected from monitoring wells MW-01, MW-03, and MW-05. The source of lead in shallow groundwater was attributed to leaching of lead to groundwater from urban fill material deposited at the site.
- Bedrock Groundwater: chlorinated VOCs (PCE and TCE)
 - OVOCs in bedrock groundwater were identified in the sample collected from monitoring well MW-13. The source of VOCs in this monitoring well may be attributed to a combination of both surficial releases of VOCs to the subsurface as well as a potential additional source from within the on-site building. The additional source has not been confirmed, and additional investigation may be warranted following the demolition of the building.
- Surface Water: No releases to surface water were identified, as indicated on Table 10.

The use and storage of chlorinated solvents at the site was discontinued when the site was vacated in 1989, thus eliminating the principal surface source of VOC contamination to site soil and groundwater.

6.3 Nature and Extent of Contamination

Based on the results of soil and groundwater monitoring conducted at the site, the lateral extent of the disposal site has been delineated and is shown on <u>Figure 2</u>. The nature and extent of contamination at the disposal site, based on the results of all Phase II site investigation activities conducted at the disposal site by Fuss & O'Neill, is detailed in <u>Section 5.0</u> and <u>Section 6.0</u>. The source of VOCs in bedrock groundwater was not completely identified, and additional investigation beneath the on-site building may be warranted.

6.4 <u>Migration Pathways</u>

A summary of the migration pathways associated with each of the types of environmental media affected by the releases at the site; sediment, soil, and groundwater, includes the following. These potential pathways are the primary methods for migration of site-related contaminants of concern (COC).

Migration and mobilization of COC in shallow soil may occur via infiltration of stormwater though vadose zone soils containing COC, and via migration of shallow groundwater through saturated soil containing COC. VOC in shallow soil can also volatilize and migrate into building structures located above soil containing VOC.



Migration of COC in shallow overburden groundwater may occur via horizontal migration though saturated overburden toward Mine Brook. Although the data presented in <u>Section 6.1</u>, indicates that the vertical gradient in the central portion of the site is generally upward, the increasing concentration of dissolved VOC in with depth indicates a migration of dissolved phase chlorinated VOC downward from shallow overburden groundwater to the deeper bedrock aquifer. The relative concentrations of COC in shallow and deep groundwater are indicative of a potential unidentified source within or beneath the abandoned building. Migration of COC in the deeper bedrock aquifer is considered to be via fracture flow, and may not be well defined by the equipotential contours representing shallow groundwater gradients. VOC in shallow groundwater can also volatilize and migrate into building structures located above the dissolved phase plume of VOC.

Migration of metals and PAH compounds through site soil and sediment and into on-site buildings is not expected due to the generally low mobility of metals and PAH compounds in soil and sediment.

7.0 RECOMMENDATIONS

Based upon the results of this assessment, the following response actions are recommended:

- On May 10, 2007 a release notification form was received by the MADEP in response to the initial Phase II ESA results. MADEP assigned RTN 2-16694 to the site. Based on MCP timelines, either a Tier Classification or Response Action Outcome will be required to be submitted to MADEP by May 10, 2008.
 - Based on the evaluation of site date relative to MCP risk-based standards, it is not feasible to conclude that a condition of "no significant risk" currently exists at the site, and therefore a permanent solution has not yet been achieved.
- The extent of the compounds of concern in environmental media at the site has been further evaluated. However, further assessment of soil, groundwater, and sediment may be required to fully delineate the releases at the site in support of MCP-related response actions and to prepare remediation plans in support of potential site redevelopment. Further assessment should be conducted beneath the on-site Lot 27 building to evaluate for additional release mechanisms and potential source areas of VOCs to bedrock groundwater.



8.0 REFERENCES

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9.0 LIMITATIONS OF WORK PRODUCT

This document was prepared for the sole use of the County of Norfolk the only intended beneficiaries of our work. Those who may use or rely upon the report and the services (hereafter "work product") performed by Fuss & O'Neill, Inc. and/or its subsidiaries or independent professional associates, subconsultants and subcontractors (collectively the "Consultant") expressly accept the work product upon the following specific conditions.

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- 2. The work product may contain information that is time sensitive. The work product was prepared by Consultant subject to the particular scope limitations, budgetary and time constraints and business objectives of the Client which are detailed therein or in the contract between Consultant and Client. Changes in use, tenants, work practices, storage, Federal, state or local laws, rules or regulations may affect the work product.
- 3. The observations described and upon which the work product was based were made under the conditions stated therein. Any conclusions presented in the work product were based solely upon the services described therein, and not on scientific or engineering tasks or procedures beyond the scope of described services.
- 4. In preparing its work product, Consultant may have relied on certain information provided by state and local officials and information and representations made by other parties referenced therein, and on information contained in the files of state and/or local agencies made available at the time of the project. To the extent that such files which may affect the conclusions of the work product are missing, incomplete, inaccurate or not provided, Consultant is not responsible. Although there may have been some degree of overlap in the information provided by these various sources, Consultant did not attempt to independently verify the accuracy or completeness of all information reviewed or received during the course of this project. Consultant assumes no responsibility or liability to discover or determine any defects in such information which could result in failure to identify contamination or other defect in, at or near the site. Unless specifically stated in the work product, Consultant assumes no responsibility or liability for the accuracy of drawings and reports obtained, received or reviewed.
- 5. If the purpose of this project was to assess the physical characteristics of the subject site with respect to the presence in the environment of hazardous substances, waste or petroleum and chemical products and wastes as defined in the work product, unless otherwise noted, no specific attempt was made to check the compliance of present or past owners or operators of the subject site with Federal, state, or local laws and regulations, environmental or otherwise.



- 6. If water level readings have been made, these observations were made at the times and under the conditions stated in the report. However, it must be noted that fluctuations in water levels may occur due to variations in rainfall, passage of time and other factors and such fluctuations may effect the conclusions and recommendations presented herein.
- 7. Except as noted in the work product, no quantitative laboratory testing was performed as part of the project. Where such analyses have been conducted by an outside laboratory, Consultant has relied upon the data provided, and unless otherwise described in the work product has not conducted an independent evaluation of the reliability of these tests.
- 8. If the conclusions and recommendations contained in the work product are based, in part, upon various types of chemical data, then the conclusions and recommendations are contingent upon the validity of such data. These data (if obtained) have been reviewed and interpretations made by Consultant. If indicated in the work product, some of these data may be preliminary or screening-level data and should be confirmed with quantitative analyses if more specific information is necessary. Moreover, it should be noted that variations in the types and concentrations of contaminants and variations in their flow paths may occur due to seasonal water table fluctuations, past disposal practices, the passage of time and other factors.
- 9. Chemical analyses may have been performed for specific parameters during the course of this project, as described in the work product. However, it should be noted that additional chemical constituents not included in the analyses conducted for the project may be present in soil, groundwater, surface water, sediments or building materials at the subject site.
- 10. Ownership and property interests of all documents, including reports, electronic media, drawings and specifications, prepared or furnished by Consultant pursuant to this project are subject to the terms and conditions specified in the contract between the Consultant and Client, whether or not the project is completed.
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- 13. Any use of or reliance on the work product shall constitute acceptance of the terms hereof.

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TABLES

FORMER NU-STYLE COMPANY, INC. PHASE II ESA ADDENDUM



Table 4 Summary of Sediment Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for the County of Norfolk, Massachusetts

February 2008

	11 1			T						
Sample Location		SD-01	SD-02	SD-03	SD-03	SD-04	SD-05*	SD-05*	SD-06*	SD-07*
Sample ID		841070426-06	841070426-07	841070426-08	841070426-09	841070426-10	937071025-01	937071025-02	937071025-04	937071025-03
Date Collected		4/26/2007	4/26/2007	4/26/2007	4/26/2007	4/26/2007	10/25/2007	10/25/2007	10/25/2007	10/25/2007
Sample Type	UNITS	Primary	Primary	Primary	Duplicate 1	Primary	Primary	Duplicate 1	Primary	Primary
Starting Depth	feet	0	0	0	0	0	0	0	0	0
Ending Depth	feet	0.2	0.2	0.2	0.2	0.2	0.25	0.25	0.25	0.25
Metals (EPA Method 6010)										
Arsenic	mg/kg	ND < 0.31	0.75	ND < 0.30	1.2	ND < 0.30	NA	NA	NA	NA
Barium	mg/kg	15	20	22	16	9.0	NA	NA	NA	NA
Beryllium	mg/kg	0.13	0.16	0.16	0.15	0.15	NA	NA	NA	NA
Cadmium	mg/kg	0.14	0.13	0.16	0.14	0.14	NA	NA	NA	NA
Chromium	mg/kg	1.6	1.3	0.75	1.1	2.9	NA	NA	NA	NA
Copper	mg/kg	7.0	6.1	1.8	1.9	3.2	NA	NA	NA	NA
Lead	mg/kg	8.6	5.9	4.8	6.6	13	NA	NA	NA	NA
Nickel	mg/kg	5.4	3.6	0.69	1.5	1.4	NA	NA	NA	NA
Thallium	mg/kg	0.55	0.99	0.69	ND < 0.30	ND < 0.30	NA	NA	NA	NA
Zinc	mg/kg	23	18	15	16	12	NA	NA	NA	NA
VOC (EPA Method 8260)										
Acetone	μg/kg	ND < 5.2	ND < 5.4	ND < 4.6	ND < 5.1	7.8	NA	NA	NA	NA
Tetrachloroethylene	μg/kg	7.6	37	ND < 4.6	ND < 5.1	ND < 4.5	NA	NA	NA	NA
Trichloroethylene	μg/kg	ND < 5.2	12	ND < 4.6	ND < 5.1	ND < 4.5	NA	NA	NA	NA
EPH/SVOC (MADEP Method/EPA Method 8270)										
C11-C22 Aromatics	μg/kg	14,000	ND < 12,000	ND < 12,000	ND < 12,000	ND < 12,000	NA	NA	NA	NA
C19-C36 Aliphatics	μg/kg	20,000	ND < 12,000	ND < 12,000	ND < 12,000	ND < 12,000	NA	NA	NA	NA
Acenaphthylene	μg/kg	140	ND < 120	ND < 120	ND < 120	ND < 120	77	160	76	170
Anthracene	μg/kg	160	ND < 120	ND < 120	ND < 120	ND < 120	340	280	58	98
Benzo(a)anthracene	μg/kg	330	ND < 120	ND < 120	ND < 120	ND < 120	440	ND < 40	400	920
Benzo(a)pyrene	μg/kg	ND < 120	ND < 120	ND < 120	ND < 120	ND < 120	330	770	350	860
Benzo(b)fluoranthene	μg/kg	120	ND < 120	ND < 120	ND < 120	ND < 120	470	690	520	1,200
Benzo(k)fluoranthene	μg/kg	140	ND < 120	ND < 120	ND < 120	ND < 120	390	1,000	450	990
Chrysene	μg/kg	ND < 120	ND < 120	ND < 120	ND < 120	ND < 120	480	1,100	500	1,200
Dibenzo(a,h)anthracene	μg/kg	ND < 120	ND < 120	ND < 120	ND < 120	ND < 120	46	100	ND < 40	160
Fluoranthene	μg/kg	820	ND < 120	ND < 120	ND < 120	ND < 120	890	2,500	810	2,000
Fluorene	μg/kg	ND < 120	ND < 120	ND < 120	ND < 120	ND < 120	ND < 38	81	ND < 40	ND < 46
Indeno (1,2,3-cd)pyrene	μg/kg	ND < 120	ND < 120	ND < 120	ND < 120	ND < 120	110	240	130	310
Phenanthrene	μg/kg	230	ND < 120	ND < 120	ND < 120	ND < 120	ND < 38	1,300	ND < 40	ND < 46
Pyrene	μg/kg	450	ND < 120	ND < 120	ND < 120	ND < 120	740	2,100	950	2,000

NOTES:

--- Not applicable ND <X: Compound not detected above laboratory reporting limit

NA: Not analyzed

NE: Not established

VOC: Volatile organic compounds

EPH: Extractable Petroleum Hydrocarbons

SVOC: Semivolatile organic compounds

* Samples collected from SD-05 through SD-07 analyzed for SVOC by EPA Method 8270; other samples analyzed for EPH by MADEP Method

Bold and color-shaded values indicate exceedance of Sediment Screening Criteria listed in Table 11.

Created by: SAH Reviewed by: TJC



Table 6 Summary of Soil Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for the County of Norfolk, Massachusetts

February 2008

Sample Location		B-02	B-02	B-04	B-04	B-05	B-05	B-06	B-06	B-10	B-10	B-11	B-11	B-12	B-15	MW-01	MW-01	MW-02	MW-02	MW-03	MW-03	MW-04	MW-04	MW-04	MW-05	MW-05	MW-13	MW-17	MW-17	MW-17
Sample ID		841061130-03	841061130-04	841061130-07	841061130-08	841061130-09	841061130-10	841061130-11	841061130-12	841061201-21	841061201-22	841061201-23	841061201-24	841061201-25	841071101-03	841061130-01	841061130-02	841061130-05	841061130-06	841061130-13	841061130-14	841061130-15	841061130-16	841061130-17	841061201-19	841061201-20	841071031-01	841071101-04	841071101-05	841071101-06
Date Collected		11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	12/1/2006	12/1/2006	12/1/2006	12/1/2006	12/1/2006	11/1/2007	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	11/30/2006	12/1/2006	12/1/2006	10/31/2007	11/1/2007	11/1/2007	11/1/2007
Sample Type	UNITS	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Primary	Duplicate 1	Primary	Primary	Primary	Primary	Primary	Primary	Duplicate 1
Starting Depth	feet	0	5	0	5	0.4	5	0	1	0	5	0	5	0	2	0.5	3	0	5	0	5	0	0	5	0.4	5	10	0.3	6	6
Ending Depth	feet	2	7	2	6	2	7	0.5	2	2	7	2	7	2	4	2	5	3	7	2	7	2	2	7	2	7	12	2	8	8
Metals (EPA Method 6010)					Ì																							Ì		
Antimony	mg/kg	ND < 0.56	ND < 0.59	ND < 0.57	ND < 0.57	ND < 0.53	ND < 0.54	ND < 0.54	ND < 0.59	ND < 0.50	ND < 0.50	ND < 0.54	ND < 0.60	ND < 0.56	NA	ND < 0.53	ND < 0.56	ND < 0.59	ND < 0.57	ND < 0.55	ND < 0.55	ND < 0.56	ND < 0.53	ND < 0.55	6.5	6.9	NA	1.0	ND < 0.18	ND < 0.18
Arsenic	mg/kg	1.2	ND < 0.59	ND < 0.57	2.0	1.1	ND < 0.54	ND < 0.54	1.8	ND < 0.50	ND < 0.50	1.8	ND < 0.60	3.1	NA	ND < 0.50	ND < 0.56	6.6	2.6	ND < 0.55	ND < 0.55	ND < 0.56	ND < 0.53	ND < 0.55	3.1	ND < 0.50	NA	3.0	1.3	0.84
Barium	mg/kg	36	20	48	24	39	10	28	26	16	9.6	23	34	30	NA	16	24	36	36	18	11	14	9.2	17	110	55	NA	26	15	15
Beryllium	mg/kg	0.21	0.19	0.34	0.36	0.25	0.13	0.18	0.16	0.7	0.91	0.16	0.38	0.26	NA	0.19	0.57	0.22	0.15	0.12	0.17	0.24	0.081	0.15	0.37	0.17	NA	0.35	0.24	0.21
Cadmium	mg/kg	0.17	0.14	0.34	0.19	0.22	ND < 0.11	0.46	0.13	ND < 0.10	ND < 0.10	0.26	ND < 0.12	0.19	NA	ND < 0.10	ND < 0.11	0.13	ND < 0.11	0.16	ND < 0.11	0.15	ND < 0.11	ND < 0.11	0.54	0.18	NA	0.6	0.37	0.17
Chromium	mg/kg	7.1	6.0	8.4	5.4	5.1	3.5	5.8	7.4	5.2	1.9	5.4	4.4	6	NA	3.2	5.3	35	4.1	2.2	5.5	6.0	1.4	5.7	27	26	NA	24	15	3.8
Copper	mg/kg	91	43	13	18	32	3.4	31	20	6.3	1.9	8.5	2.9	37	NA	4.9	12	160	9.0	5.0	2.9	2.0	2.5	25	29	9.5	NA	110	11	3.1
Lead	mg/kg	40	18	8.4	22	20	1.6	97	25	2.9	4.8	17	4.3	93	NA	4.9	8.1	25	89	9.2	2.6	3.4	1.5	4.7	780	310	NA	68	7.1	2.5
Mercury (EPA Method 7471)	mg/kg	0.029	$ND \le 0.024$	0.034	0.051	0.023	$ND \le 0.022$	ND < 0.021	0.065	0.023	ND < 0.021	0.032	ND < 0.024	0.044	NA	ND < 0.021	ND < 0.022	0.14	ND < 0.023	ND < 0.022	ND < 0.022	ND < 0.023	ND < 0.021	ND < 0.022	0.073	ND < 0.023	NA	0.12	0.028	ND < 0.024
Nickel	mg/kg	4.0	3.6	23	37	4.9	14	10	2.6	3.6	1.0	3.2	1.7	130	NA	2.6	3.3	6.2	5.0	3.2	1.8	2.0	6.5	2.0	6.4	6.3	NA	4.3	4.7	2.2
Zinc	mg/kg	85	63	20	26	48	6.8	71	14	22	15	48	8.4	28	NA	10	13	27	54	14	6.3	4.0	4.2	16	310	84	NA	73	28	11
VOC (EPA Method 8260)																														
1,1,1-trichloroethane	μg/kg	ND < 5.1	ND < 5.8	ND < 570	ND < 1,100	ND < 5.0	ND < 5.3	ND < 1,100	ND < 5.8	ND < 270	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	73	17	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Acetone	μg/kg	ND < 20	ND < 23	ND < 2,300	ND < 4,500	ND < 20	ND < 21	ND < 4,300	ND < 23	ND < 1,100	ND < 22	ND < 20	30	ND < 21	ND < 1,200	ND < 21	ND < 22	ND < 21	ND < 22	ND < 21	ND < 22	ND < 21	ND < 20	ND < 18	ND < 18	ND < 22	ND < 3.2	11	28	35
M/P-xylenes	μg/kg	7.0	ND < 5.8	ND < 350	ND < 340	ND < 5.0	ND < 5.3	ND < 300	ND < 5.8	ND < 270	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	ND < 5.2	ND < 4.9	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Methyl ethyl Ketone	μg/kg	ND < 10	ND < 12	ND < 1,100	ND < 2,300	ND < 10	ND < 10	ND < 2,100	ND < 12	ND < 550	ND < 11	ND < 10	ND < 11	ND < 10	ND < 1,200	ND < 10	ND < 11	ND < 11	ND < 11	ND < 10	ND < 11	ND < 10	ND < 9.8	ND < 9.2	ND < 9.2	ND < 11	ND < 3.2	ND < 6.4	6.6	7.4
Naphthalene	μg/kg	ND < 5.1	10	ND < 110	ND < 110	ND < 5.0	ND < 5.3	ND < 110	ND < 5.8	ND < 100	ND < 5.4	ND < 5.0	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	260	ND < 5.4	ND < 5.2	2300	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	140	ND < 4.7	ND < 5.2
Tetrachloroethylene	μg/kg	ND < 5.1	ND < 5.8	15,000	20,000	110	22	34,000	310	4,300	48	40	45	11	40,000	ND < 5.3	ND < 5.6	28	45	130	120	13	18	26	ND < 4.6	ND < 5.6	6.4	ND < 6.4	ND < 4.7	ND < 5.2
Toluene	μg/kg	17	ND < 5.8	ND < 350	ND < 340	ND < 5.0	ND < 5.3	ND < 300	ND < 5.8	ND < 270	ND < 5.4	16	ND < 5.7	ND < 5.2	ND < 1,200	ND < 5.3	ND < 5.6	ND < 5.3	ND < 5.5	ND < 5.2	ND < 5.4	ND < 5.2	ND < 4.9	ND < 4.6	ND < 4.6	ND < 5.6	ND < 3.2	ND < 6.4	ND < 4.7	ND < 5.2
Trichloroethylene	μg/kg	ND < 5.1	ND < 5.8	19,000	31,000	58	9.6	6,700	79	9,300	150	5	ND < 5.7	6.5	9,200	ND < 5.3	ND < 5.6	12	21	150	67	37	44	24	ND < 4.6	ND < 5.6	3.5	ND < 6.4	ND < 4.7	ND < 5.2
EPH with Targets (MADEP Method)																														
C11-C22 Aromatics	μg/kg	17,000	32,000	40,000	61,000	33,000	ND < 11,000	92,000	ND < 12,000	25,000	ND < 10,000	18,000	ND < 11,000	28,000	NA	100,000	16,000	110,000	20,000	52,000	81,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	ND < 11,000	NA	60,000	17,000	ND < 12,000
C19-C36 Aliphatics	μg/kg	ND < 11,000	ND < 11,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	24,000	ND < 12,000	ND < 10,000	ND < 10,000	ND < 11,000	ND < 11,000	16,000	NA	40,000	14,000	ND < 11,000	ND < 11,000	ND < 10,000	38,000	ND < 11,000	ND < 11,000	ND < 10,000	ND < 11,000	ND < 11,000	NA	18,000	ND < 12,000	ND < 12,000
2-Methylnaphthalene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	210	ND < 110	200	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	ND < 130	ND < 120	ND < 120
Acenaphthene	μg/kg	740	ND < 110	ND < 110	ND < 110	220	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	ND < 110	ND < 110	560	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	150	ND < 120	ND < 120
Acenaphthylene	μg/kg	240	340	230	250	ND < 100	ND < 110	260	ND < 120	200	ND < 100	ND < 110	ND < 110	120	NA	270	ND < 100	1,300	350	150	120	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	260	ND < 120	ND < 120
Anthracene	μg/kg	ND < 110	320	200	ND < 110	970	ND < 110	340	ND < 120	250	ND < 100	ND < 110	ND < 110	ND < 110	NA	390	ND < 100	2,000	ND < 110	1,800	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	640	ND < 120	ND < 120
Benzo(a)anthracene	μg/kg	ND < 110	1,000	330	310	1,100	ND < 110	560	ND < 120	1,500	ND < 100	ND < 110	ND < 110	490	NA	1,200	ND < 100	4,400	200	2,800	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,800	ND < 120	ND < 120
Benzo(a)pyrene	μg/kg	ND < 110	1,000	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	2,000	ND < 100	ND < 110	ND < 110	ND < 110	NA	1,100	ND < 100	3,900	ND < 110	2,200	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,600	ND < 120	ND < 120
Benzo(b)fluoranthene	μg/kg	ND < 110	1,400	560	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	2,000	ND < 100	ND < 110	ND < 110	290	NA	1,700	ND < 100	5,600	ND < 110	2,600	230	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,800	ND < 120	ND < 120
Benzo(k)fluoranthene	μg/kg	ND < 110	540	170	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	860	ND < 100	ND < 110	ND < 110	ND < 110	NA	600	ND < 100	1,900	ND < 110	1,100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,300	ND < 120	ND < 120
Chrysene	μg/kg	ND < 110	1,200	640	590	1,200	ND < 110	810	ND < 120	120	ND < 100	ND < 110	ND < 110	ND < 110	NA	1,600	ND < 100	4,500	280	2,800	440	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,700	ND < 120	ND < 120
Fluoranthene	μg/kg	490	2,100	380	410	2,300	ND < 110	940	ND < 120	1,300	ND < 100	ND < 110	ND < 110	1,300	NA	2,200	ND < 100	8,200	320	6,700	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	3,900	180	ND < 120
Fluorene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	230	ND < 110	ND < 110	ND < 120	ND < 100	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	390	ND < 110	630	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	280	ND < 120	ND < 120
Indeno (1,2,3-cd)pyrene	μg/kg	ND < 110	ND < 110	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	ND < 120	260	ND < 100	ND < 110	ND < 110	ND < 110	NA	ND < 100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	1,200	ND < 120	ND < 120
Phenanthrene	μg/kg	350	1,100	150	190	2,300	ND < 110	470	ND < 120	260	ND < 100	ND < 110	ND < 110	940	NA	990	ND < 100	3,500	200	7,600	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	2,800	290	ND < 120
Pyrene	μg/kg	ND < 110	2,000	470	410	2,000	ND < 110	990	ND < 120	1,600	ND < 100	ND < 110	ND < 110	1,200	NA	2,200	ND < 100	6,200	310	6,000	ND < 100	ND < 110	ND < 110	ND < 100	ND < 110	ND < 110	NA	3,600	180	ND < 120
VPH with Targets (MADEP Method)																-														
C9-C12 Aliphatics	μg/kg	ND < 6,900	ND < 8,800	ND < 7,000	ND < 6,900	ND < 5,900	ND < 6,500	ND < 5,900	ND < 7,800	ND < 6,400	ND < 6,100	ND < 12,000	ND < 8,000	ND < 6,800	NA	ND < 6,200	ND < 6,500	ND < 7,900	ND < 7,000	ND < 6,400	ND < 6,400	ND < 7,600	ND < 6,000	ND < 6,100	14,000	ND < 7,600	NA	ND < 6,400	ND < 6,100	ND < 6,000

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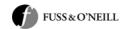


Table 8 Summary of Groundwater Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for the County of Norfolk, Massachusetts

February 2008

Sample Location		MCP Regulat	ory Standards	MW-01	MW-01	MW-01	MW-02	MW-02	MW-03	MW-03	MW-04	MW-05	MW-05	MW-13	MW-13	MW-14	MW-16	MW-17	MW-17
Sample ID		MA Method 1 GW	MA Method 1 GW	841061208-27	841061208-28	841071106-03	841061208-30	841071107-10	841061208-32	841071107-11	841061208-29	841061208-31	841071107-09	841071106-05	841071106-06	841071106-04	841071107-08	841071106-01	841071106-02
Date Collected		Std Application	Std Application	12/8/2006	12/8/2006	11/6/2007	12/8/2006	11/7/2007	12/8/2006	11/7/2007	12/8/2006	12/8/2006	11/7/2007	11/6/2007	11/6/2007	11/6/2007	11/7/2007	11/6/2007	11/6/2007
Sample Type		for GW-2 area	for GW-3 area	Primary	Duplicate	Primary	Duplicate	Primary	Primary	Primary	Duplicate								
Groundwater Parameters	UNITS																		
рН	SU			5.97	5.97	6.35	6.59	6.86	6.33	5.91	5.97	NA	6.03	5.68	5.68	5.78	5.91	6.36	6.36
Specific Conductance	μS/cm			464	464	470	1,727	1,547	1,534	3,129	2,010	NA	NA	3,761	3,761	3,065	2,543	494	494
Temperature	C deg			13.1	13.1	17	8.3	13.7	7.6	14.1	10.5	NA	13.2	12	12	15.5	14.9	17.9	17.9
Turbidity	ntu			36	36	20.5	50	NA	500	NA	15	NA	NA	4.64	4.64	3.3	8.9	4.4	4.4
Dissolved Oxygen	mg/l			0.4	0.4	0.1	7.7	4.9	6.9	4.9	2.6	NA	NA	2.8	2.8	3.5	1.6	0.1	0.1
ORP	mv			-35.0	NA	-116.2	59	45.8	93.2	51.3	45.1	NA	NA	30	30	20.2	52.6	-53.2	NA
Metals (EPA Method 6010)																			
Barium, Total	mg/l	NE	50	0.042	0.038	0.031	0.15	NA	0.21	NA	0.14	0.83	NA	0.24	0.25	0.12	0.11	0.06	0.061
Barium, Dissolved	mg/l	NE	50	NA	NA	NA	NA	0.18	NA	0.17	NA	NA	0.39	NA	NA	NA	NA	NA	NA
Beryllium	mg/l	NE	0.05	ND < 0.0010	ND < 0.0010	NA	ND < 0.0010	NA	0.0087	NA	ND < 0.0010	0.0018	NA						
Cadmium	mg/l	NE	0.004	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	NA	ND < 0.0020	NA	ND < 0.0020	0.0034	NA	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020
Chromium	mg/l	NE	0.3	ND < 0.010	ND < 0.010	0.0029	ND < 0.010	NA	0.036	NA	ND < 0.01	0.092	NA	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020
Copper	mg/l	NE	NE	ND < 0.010	ND < 0.010	NA	0.015	NA	0.018	NA	ND < 0.01	0.073	NA						
Lead, Total	mg/l	NE	0.01	0.014	0.012	0.0066	ND < 0.0040	NA	0.098	NA	ND < 0.0040	1.9	NA	0.0033	0.0053	ND < 0.0020	ND < 0.0020	ND < 0.0020	ND < 0.0020
Lead, Dissolved	mg/l	NE	0.01	NA	NA	NA	NA	0.0026	NA	0.006	NA	NA	0.094	NA	NA	NA	NA	NA	NA
Nickel	mg/l	NE	0.2	ND < 0.010	ND < 0.010	NA	0.15	NA	0.054	NA	0.017	0.12	NA						
Zinc	mg/l	NE	0.9	0.023	0.015	NA	0.057	NA	0.17	NA	0.028	0.73	NA						
VOC (EPA Method 8260)																			
1,1,1-trichloroethane	μg/l	4,000	20,000	ND < 1.0	1.8	ND < 1.0													
cis-1,2-dichloroethylene	μg/l	100	50,000	ND < 1.0	ND < 1.0	ND < 1.0	ND < 1.0	1.7	ND < 1.0	ND < 1.0	8	ND < 1.0							
Methyl tert butyl ether	μg/l	50,000	50,000	ND < 1.0	1.8	ND < 1.0	ND < 1.0	1.5	1.4	ND < 1.0	ND < 1.0	ND < 1.0	ND < 1.0						
Tetrachloroethylene	μg/l	50	30,000	ND < 1.0	ND < 1.0	ND < 1.0	6.6	23	43	74	240	ND < 1.0	1.3	290	260	12	41	ND < 1.0	ND < 1.0
Trichloroethylene	μg/l	30	5,000	ND < 1.0	ND < 1.0	ND < 1.0	6.6	25	40	59	150	ND < 1.0	ND < 1.0	60	56	20	45	ND < 1.0	ND < 1.0

NOTES:

--- Not applicable
ND <X: Compound not detected above laboratory reporting limit

NA: Not analyzed

NE: Not established

μS/cm: microsiemens per centimeter

C deg: degrees Celcius

ntu: nephelometric turbidity units ORP: Oxidation-reduction potential

mv: millivolts

mg/l: milligrams per liter μg/l: micrograms per liter

VOC: Volatile organic compounds Bold and color-shaded values indicate exceedence of one or more regulatory criteria.

Created by: SAH Reviewed by: TJC



Table 10 Summary of Surface Water Analytical Data and Objectives

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for the County of Norfolk, Massachusetts

February 2008

				CWI 04	OWI OO	CWI OZ	OWI OZ	CWI O.4
				SW-01	SW-02	SW-03	SW-03	SW-04
		MA Method 1 GW	US EPA Chronic	841070426-01	841070426-02	841070426-03	841070426-04	841070426-05
		Std Application	Criteria Continuous	04/26/2007	04/26/2007	04/26/2007	04/26/2007	04/26/2007
		for GW-3 area	Concentrations	Primary	Primary	Primary	Duplicate	Primary
Metals (Method 6010)	·		_		_			
Barium	mg/l	50	NE	0.086	0.085	0.084	0.083	0.083
Copper	mg/l	NE	0.0090	0.0040	0.0023	ND < 0.0020	0.0041	0.0023
Lead	mg/l	0.01	0.0025	ND < 0.0020	ND < 0.0020	0.0033	ND < 0.0020	ND < 0.0020
Zinc	mg/l	0.9	0.12	0.018	0.017	0.017	0.016	0.015
VOC (Method 8260)		Varies	Varies	ND < varies	ND < varies	ND < varies	ND < varies	ND < varies
VPH (MADEP Method)								
Methyl tert-butyl ether (MTBE)	μg/l	50,000	NE	ND < 1.0	1.1	ND < 1.0	1.1	ND < 1.0
EPH (MADEP Method)	μg/1	Varies	Varies	ND < varies	ND < varies	ND < varies	ND < varies	ND < varies

Created by: SAH

Reviewed by: TJC

ND <X: Compound not detected above laboratory reporting limit

US EPA: United States Environmental Protection Agency

VPH: Volatile petroleum hydrocarbons EPH: Extractable petroleum hydrocarbons

VOC: Volatile organic compounds

NE: Not established



Table 11 Summary of MADEP Criteria for Detected Compounds in Soil and Sediment

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for the County of Norfolk, Massachusetts

February 2008

		Regulatory Criteria for Soil						
		MA Method 1 S-1	MA Method 1 S-1	MA Method 1 S-2	MA Method 1 S-2	MA Method 1 S-3	MA Method 1 S-3	MADEP Stage I
		Std Application	Std Application	Std Application	Std Application	Std Application	Std Application	Freshwater Sediment
		for GW-2 area	for GW-3 area	for GW-2 area	for GW-3 area	for GW-2 area	for GW-3 area	Screening Criteria*
Total Metals (via Method 6010/7471)	UNITS		•					
Antimony	mg/kg	20	20	30	30	30	30	NE
Arsenic	mg/kg	20	20	20	20	20	20	33
Barium	mg/kg	1,000	1,000	3,000	3,000	5,000	5,000	NE
Beryllium	mg/kg	1	1	1	1	3	3	NE
Cadmium	mg/kg	2	2	30	30	30	30	5.0
Chromium	mg/kg	30	30	200	200	200	200	110
Copper	mg/kg							150
Lead	mg/kg	300	300	300	300	300	300	130
Mercury	mg/kg	20	20	30	30	30	30	0.18
Nickel	mg/kg	20	20	700	700	700	700	49
Thallium	mg/kg	8,000	8,000	60,000	60,000	80,000	80,000	NE
Zinc	mg/kg	2,500	2,500	3,000	3,000	5,000	5,000	460
VPH (MADEP Method)								
C9-C12 Aliphatics	μg/kg	1,000,000	1,000,000	2,500,000	2,500,000	5,000,000	5,000,000	
VOC (Method 8260B)								
1,1,1-trichloroethane	μg/kg	500,000	500,000	600,000	1,000,000	600,000	3,000,000	NE
Acetone	μg/kg	60,000	60,000	60,000	60,000	60,000	60,000	NE
M/P-xylenes	μg/kg	300,000	300,000	5,000,000	5,000,000	300,000	300,000	NE
Methyl ethyl Ketone	μg/kg	40,000	40,000	40,000	40,000	40,000	40,000	NE
Naphthalene	μg/kg	40,000	500,000	2,000,000	2,000,000	40,000	3,000,000	NE
Tetrachloroethylene	μg/kg	10,000	30,000	10,000	200,000	10,000	1,000,000	NE
Toluene	μg/kg	300,000	500,000	2,500,000	2,500,000	300,000	1,000,000	NE
Trichloroethylene	μg/kg	2,000	90,000	2,000	700,000	2,000	2,000,000	NE
EPH (MADEP Method)								
C19-C36 Aliphatics	μg/kg	2,500,000	2,500,000	5,000,000	5,000,000	5,000,000	5,000,000	NE
C11-C22 Aromatics	μg/kg	800,000	800,000	2,000,000	2,000,000	5,000,000	5,000,000	NE
2-Methylnaphthalene	μg/kg	500,000	500,000	1,000,000	1,000,000	2,000,000	1,000,000	NE
Acenaphthene	μg/kg	1,000,000	1,000,000	2,500,000	2,500,000	5,000,000	4,000,000	NE
Acenaphthylene	μg/kg	100,000	100,000	2,500,000	1,000,000	2,500,000	1,000,000	NE
Anthracene	μg/kg	1,000,000	1,000,000	2,500,000	2,500,000	5,000,000	5,000,000	57
Benzo(a)anthracene	μg/kg	7,000	7,000	40,000	40,000	300,000	300,000	110
Benzo(a)pyrene	μg/kg	2,000	2,000	4,000	4,000	30,000	30,000	150
Benzo(b)fluoranthene	μg/kg	7,000	7,000	40,000	40,000	300,000	300,000	NE
Benzo(ghi)perylene	μg/kg	1,000,000	1,000,000	2,500,000	2,500,000	2,500,000	2,500,000	NE
Benzo(k)fluoranthene	μg/kg	70,000	70,000	400,000	400,000	3,000,000	3,000,000	NE
Chrysene	μg/kg	7,000	7,000	10,000	10,000	40,000	40,000	170
Dibenzo(a,h)anthracene	μg/kg	700	700	4,000	4,000	30,000	30,000	33
Fluoranthene	μg/kg	1,000,000	1,000,000	3,000,000	3,000,000	5,000,000	5,000,000	420
Fluorene	μg/kg	1,000,000	1,000,000	3,000,000	2,000,000	5,000,000	4,000,000	77
Indeno (1,2,3-cd)pyrene	μg/kg	7,000	7,000	40,000	40,000	300,000	300,000	NE
Naphthalene	μg/kg	40,000	500,000	2,000,000	2,000,000	40,000	3,000,000	180
Phenanthrene	μg/kg	1,000,000	100,000	2,500,000	100,000	2,500,000	100,000	200
Pyrene	μg/kg	1,000,000	1,000,000	3,000,000	3,000,000	5,000,000	5,000,000	200

NOTES:

µg/kg: micrograms per kilogram VOC: volatile organic compounds mg/kg: milligrams per kilogram EPH: Extractable Petroleum Hydrocarbons

NE: not established MADEP: Massachusetts Department of Environmental Protection
S: soil *Sediment Screening Criteria incorporate Threshold Effect Concentrations
GW: groundwater of MacDonald et al. (2000) and revised Sediment Screening Criteria published

VPH: Volatile Petroleum Hydrocarbons by MADEP (2005).

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Table 12 Groundwater Elevation Measurements for On-Site Monitoring Wells Gauged November 7, 2007

Former Nu-Style Property Grove Street Franklin, Massachusetts

Phase II Environmental Site Assessment Addendum Report Prepared for Norfolk County, Massachusetts

February 2008

Location	Time	Depth to Water (feet from PVC)	Absolute Elevation of PVC ^a (feet)	Groundwater Elevation (feet)		
MW-1 ^b	0930	4.60	100.35	95.75		
MW-2	0911	8.16	98.54	90.38		
MW-3	0920	8.07	99.73	91.66		
MW-4		U	nable to locate MW-4			
MW-5	1100	9.20	104.47	95.27		
MW-13	0915	6.99	99.31	92.32		
MW-14	0926	10.41	104.40	93.99		
MW-16	0924	7.15	100.81	93.66		
MW-17 ^b	0920	4.61	100.37	95.76		

^aelevation data from surveys conducted December 4, 2006 and November 7, 2007

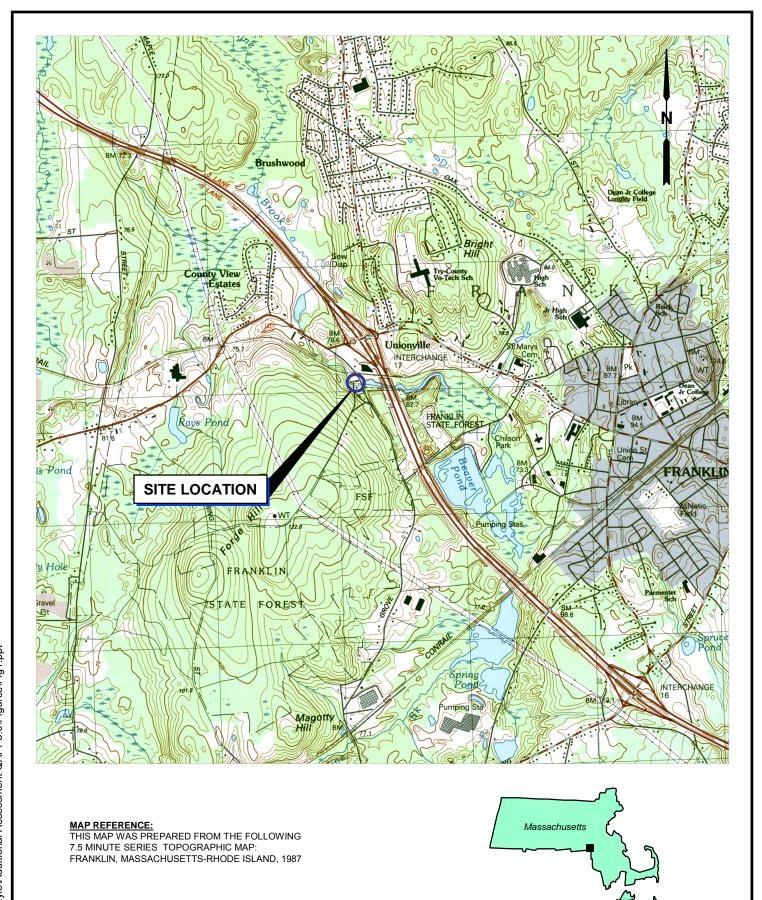
Created by SAH Reviewed by TJC

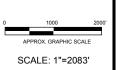
^bMonitoring wells MW-1 and MW-17 were gauged on November 6, 2007 survey data utilize the same arbitrary 100.00-foot benchmark



FIGURES

FORMER NU-STYLE PROPERTY PHASE II ESA ADDENDUM







FOUNDRY CORPORATE OFFICE CENTER
275 PROMENADE ST, SUITE 350, PROVIDENCE RI 02908
401-861-3070 www.FandO.c

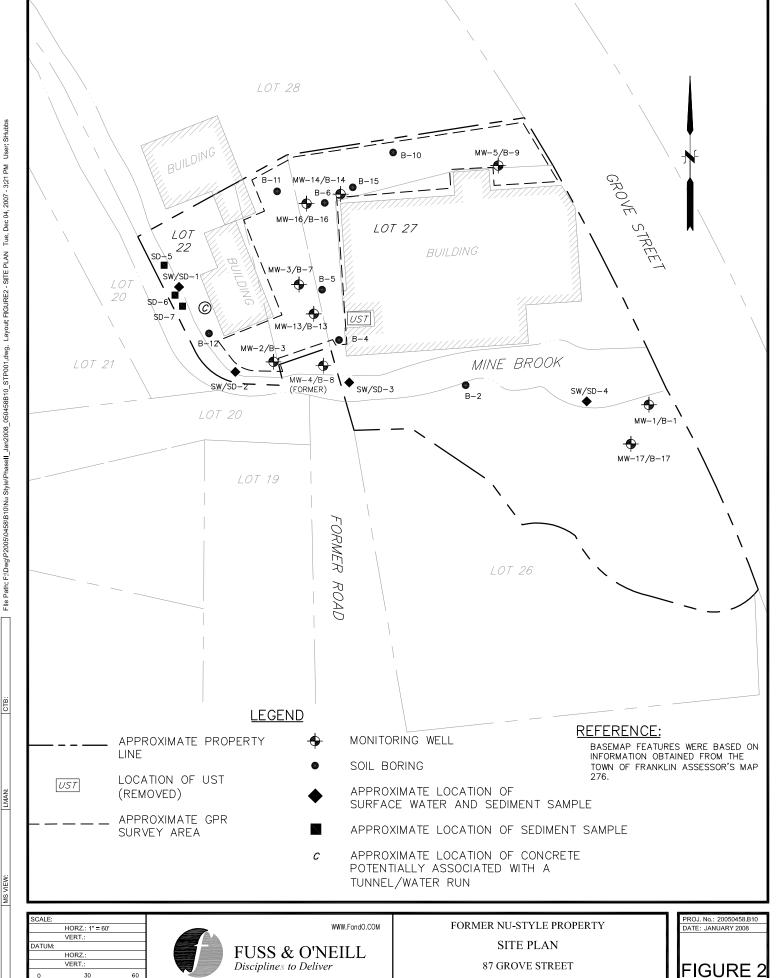
NORFOLK COUNTY, MASSACHUSETTS
SITE LOCATION MAP

FORMER NU-STYLE COMPANY, INC.

87 GROVE STREET FRANKLIN, MASS.

PROJ. No: 20050458.B10 DATE: JANUARY 2008

FIGURE 1



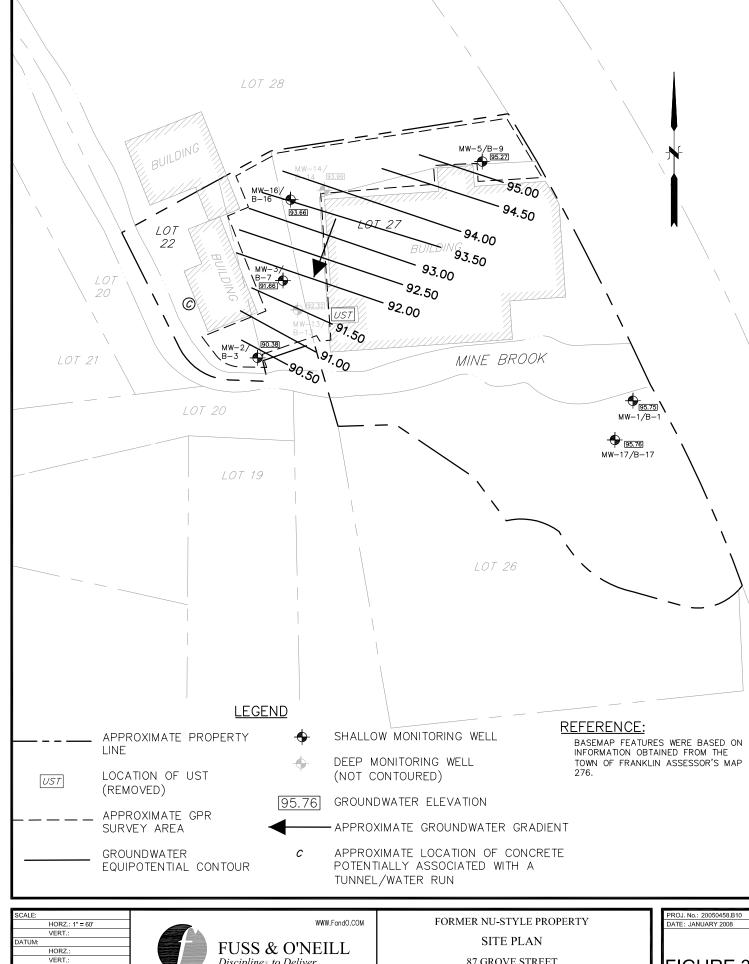
FRANKLIN

MASSACHUSETTS

275 PROMENADE ST SUITE 350 PROVIDENCE RI 02908 401.861.3070

-

GRAPHIC SCALE



GRAPHIC SCALE

LMAN:

MS VIEW:

File Path: F:Dwg/P2005/0458/B10/Nu Style/Phasell_Jan2008_050458B10_GW_CONTOUR001.dwg, Layout: FIGURE3 - GW CONTOUR MAP Tue, Dec 04, 2007 - 3:15 PM User: SHubbs

Disciplines to Deliver

275 PROMENADE ST SUITE 350 PROVIDENCE RI 02908 401.861.3070

87 GROVE STREET

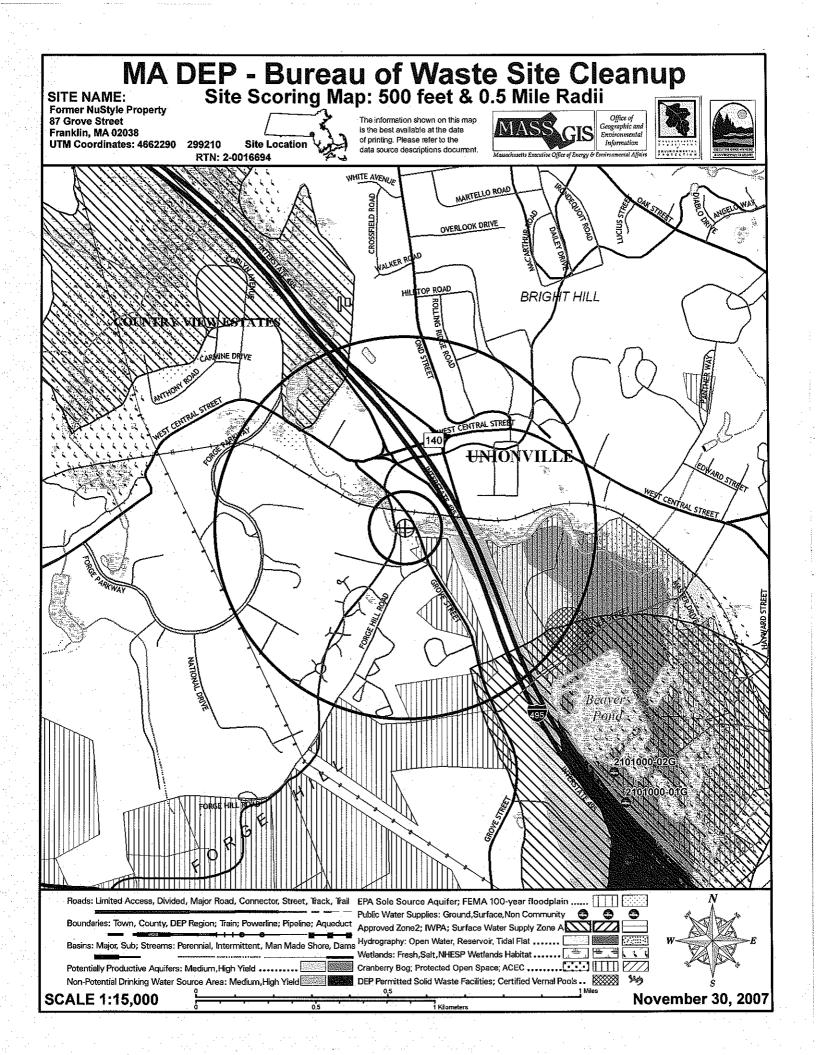
MASSACHUSETTS FRANKLIN

FIGURE 3



APPENDIX A

MADEP BUREAU OF WASTE SITE CLEANUP SITE SCORING MAP





APPENDIX B

SOIL BORING LOGS AND MONITORING WELL COMPLETION REPORTS

Township/Range: Franklin, Massachusetts

Site Id: MW-13

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Project Number: 2005-0458 B10



Location: Center of Old Grove Street Datum:

Description: Monitoring Well, Shallow

Date(s): 10/31/07 - 10/31/07

Completed Depth: 35.00'

Total Depth: 35.00'

Drilling Method: Hollow Stem Auger/Air Rotary

Remarks: Field Instrument: OVM MiniRAE

Development Method: Surge block on 11/05/2007.

No refusal.

Driller: Phil & Brian Logged By: S. Hubbs

Contractor: Subsurface

Borehole Dia.: 8 in Corehole Dia.: 3.00in

Blank Casing: type: PVC

dia: 2.00in fm: 0.0' to: 25.00'

Screens:

type: Slotted size: 0.010in dia: 2.00in fm: 25.00' to: 35.00'

Annular Fill:

type: Concrete fm: 0.00' to: 0.50' type: Sand and Native Material to: 12.50' fm: 0.50' to: 23.00' to: 35.00' type: Bentonite Pellets fm: 12.50' type: #1 Sand

fm: 23.00'

					type:		fm:		to:	
Elevation	Depth	Sample No.	Recovery	Blow Count (SPT Test)	Material Description	Graphic Log	USCS Code	Well Constru		OVM
-	-	N/A		9 8 8 9	0-0.4': SAND, F; some silt; trace asphalt; dusky brown (5YR 2/2), dry. (Fill). 0.4-2.0': Sand, F-M and silt; some gravel; moderate brown (5YR 4/4), dry.		FI SW	<u> </u>		O ppm
-2-	2 -	N/A		5 4 4 4	SAND, F-M; trace gravel; dusky yellow (5Y 6/4), dry.	· · · · · · · · · · · · · · · · · · ·	311			0.5 ppm
-4-	4	N/A		3 1 2 1	Same as above.		SP			0.9 ppm
-6- -	6	N/A	7	3 3 9 21	Sand, F-M and silt and gravel; moderate brown (5YR 4/4), dry.					O ppm
-8-	8	N/A		11 10 11 19	Same as above, wet.	0 0	SW			0.3 ppm
-10 -	10 -	O1	7	10 27 36 61	Sand, C and gravel; some F-M sand and silt; moderate olive brown (5Y 4/4), wet.	0 0 0				0.8 ppm
-12-	12	N/A		10	Same as above. BEDROCK, grantitic.					0.3 ppm
-14 -	14									
-16 - Chec	16-		K					Page	1 of	2

Township/Range: Franklin, Massachusetts

Site Id: MW-13

Project Number: 2005-0458 B10



Section Sect							275	PROMENADE	STREET, SUIT	E 350, PROVIDENCE, RI 02908	
-20 - 20	Elevation	Depth	Sample No.	Recovery	Blow Count	Material Description		Graphic Log	USCS Code	Well Construction	ОУМ
-22 - 22	-18-	18-									
GR -24 - 24 - 24 -	-20-	20 —									
-26 - 26 - 26 - 28 - Water bearing fracture. -30 - 30 - 30 - 32 - 32 - 34 - 34 - 34 - 36 - 36 - 36 - 36 - 36	-22-	22 –									
-28 - 28 - Water bearing fracture. -30 - 30	-24 -	24							GR	_	
-30 - 30	-26- -	26-									
-32 - 3234 - 3436 - 3636 - 3636 - 36 -	-28 -	28 -				Water bearing fracture.					
-34 - 34 - End of boring at 35 feet.	-30 -	30 -									
End of boring at 35 feet.	-32 <i>-</i>	32 —									
	-34 -	34 —									
_38 - 38 -	-36-	36-				End of boring at 35 feet.					
· · · · · · · · · · · · · · · · · · ·	-38-	38 —									

Township/Range: Franklin, Massachusetts

Site Id: MW-14

Ground Elevation: 0.00'

Coordinate X: 0.00

Coordinate Y: 0.00

Project Number: 2005-0458 B10



Location: NW corner lot 27 building Datum:

Description: Monitoring Well, Shallow

Date(s): 10/31/07 - 10/31/07

Completed Depth: 21.00'

Total Depth: 21.00'

Drilling Method: Hollow Stem Auger/Air Rotary

Remarks: Field Instrument: None

Development Method: Surge block on 11/05/2007.

No refusal.

Driller: Phil & Brian Logged By: S. Hubbs

Contractor: Subsurface

Borehole Dia.: 8 in Corehole Dia.: 3.00in

Blank Casing: type: PVC

fm: -2.3'dia: 2.00in to: 11.00'

Screens:

type: Slotted size: 0.010in dia: 2.00in fm: 11.00' to: 21.00'

Annular Fill:

type: Concrete fm: 0.00' to: 0.50' type: Sand and Native Material fm: 0.50' to: 8.00' type: Bentonite Pellets fm: 8.00' to: 10.00' type: #1 Sand fm: 10.00' to: 21.00'

type: fm: to:

	цуре: 		TM:	to:
Elevation Depth Sample No. Recovery Blow Count	Material Description	Graphic Log	SSS C	Vell Construction
-2- 2- -4- 4- -6- 6- -8- 8-	No samples. See boring log for B—15 for material descriptions.	Or Control of the Con	FI SM	MP. EL. O.GO
-10 - 10	BEDROCK.		SH	Page 1 of 2

Township/Range: Franklin, Massachusetts

Site Id: MW-14

Project Number: 2005-0458 B10



						275	PROMENADE	STREET, SUIT	E 350, PROVIDENCE, RI 02908	
Elevation	Depth	Sample No.	Recovery	Blow Count	Material Description		Graphic Log	USCS Code	Well Construction	Vapor
-18-	18 —									
-20-	20 —									
-22-	22 —				End of boring at 21 feet.		}1U1U		Ken Ken	
- -24 -	24 —									
-26-	26-									
-28-	28 –									
-30 -	30 —									
- -32 -	32 —									
-34 -	- 34 -									
-36-	- 36 —									
-38 <i>-</i>	- 38 —									
		<u> </u>	1		1		I		1	

Site Id: B-15

Project Number: 2005-0458 B10



Project Location: Franklin, Massachusetts Location: NW corner lot 27 building Driller: Phil & Brian Datum: Logged By: S. Hubbs Description: Soil Boring Ground Elevation: 0.00' Contractor: Subsurface Borehole Dia.: 8.00in Date(s): 11/01/07 - 11/01/07 Coordinate X: 0.00 Drilling Method: Hollow Stem Auger Total Depth: 6.00' Coordinate Y: 0.00 Back Fill: Remarks: Field Instrument: OVM MiniRAE type: Native Material fm: 0.00' to: 6.00' type: fm: to:

					type: fm: type: fm: type: fm: type: fm:		to: to:	
					type: fm: type: fm:		to: to:	
Elevation	Depth	Sample No.	Recovery	Blow Count (SPT Test)	Material Description	Graphic Log	USCS Code	MVO
-	_	N/A		3566	Sand, F-M and gravel and coal and slag; dusky brown (5YR 2/2). (Fill).		FI	9 ppm
-2-	2	-03		3 5 5 9	Sand, F and silt; trace gravel; moderate brown (5YR 4/4), moist.	X	SM	9.5 ppm
-4-	4	N/A	/	21 43 66 100	BEDROCK.		SH	
-6 -	6 —				Refusal and end of boring at 6.0 feet.			
-8-	8-							
-10-	10 —							
-12-	12 —							
- -14	14 —							
-16 <i>-</i> -	16 —							
- -18-	18 —							
Checke	ed By:	BEK			Po	age 1	of 1	

Township/Range: Franklin, Massachusetts

Site Id: MW-16

Coordinate X: 0.00

Coordinate Y: 0.00

Project Number: 2005-0458 B10



Driller: Phil & Brian

Borehole Dia.: 8.00in

Location: NW corner lot 27 building Description: Monitoring Well, Shallow

Date(s): 11/01/07 - 11/01/07 Completed Depth: 10.00'

Total Depth: 10.00'

Remarks: Field Instrument: OVM MiniRAE

Development Method: Geopump

No refusal.

Datum: Logged By: S. Hubbs Contractor: Subsurface Ground Elevation: 0.00'

Drilling Method: Hollow Stem Auger

Blank Casing: type: PVC

dia: 2.00in fm: 0.0' to: 5.00'

Screens:

type: Slotted size: 0.010in dia: 2.00in fm: 5.00' to: 10.00'

Annular Fill:

fm: 0.00' to: 0.50' type: Concrete type: Sand and Native Material fm: 0.50' to: 2.00' type: Bentonite Chips to: 3.00' fm: 2.00'

type: #1 Sand fm: 3.00' to: 10.00'

type: fm: to: Log င္ပဝရွိ ģ Count Test) Material Description Elevation Sample I Recovery Well Construction Graphic Depth MP. EL. 0.00 0-7.0': No samples. -2-2 4 SW -6 6 N/A 100 0.6 ppm Sand, F and silt and gravel; moderate olive brown (5Y 4/4), moist. -8 8 SH **-10** -10 End of boring at 10 feet. **-12** -12 -14 -14 -16 16 BEK Checked By: Page 1 of 1

Township/Range: Franklin, Massachusetts

Site Id: MW-17

Project Number: 2005-0458 B10



Location: Parking lot, S of Mine Brook Datum:

Description: Monitoring Well, Shallow

Date(s): 11/01/07 - 11/01/07

Ground Elevation: 0.00'

Coordinate X: 0.00

Completed Depth: 14.00'

Coordinate Y: 0.00

Total Depth: 14.00'

Remarks: Field Instrument: OVM MiniRAE

Development Method: Surge Block on 11/05/2007.

No refusal.

Logged By: S. Hubbs Driller: Phil & Brian Contractor: Subsurface Borehole Dia.: 8.00in

Drilling Method: Hollow Stem Auger

Blank Casing: type: PVC

dia: 2.00in fm: 0.0' to: 4.00'

Screens:

type: Slotted size: 0.010in dia: 2.00in fm: 4.00' to: 14.00'

Annular Fill:

type: Concrete fm: 0.00' to: 0.50' type: Sand and Native Material fm: 0.50' to: 1.00' type: Bentonite Chips to: 2.50' fm: 1.00' type: #1 Sand fm: 2.50' to: 14.00'

					type: #1 Sand type:		fm: fm:	2.50' to: 1 to:	4.00
Elevation	Depth	Sample No.	Recovery	Blow Count (SPT Test)	Material Description	Graphic Log	USCS Code	Well Construction MP. EL. 0.00	OVM
-	_	N/A -04		3 7 5 4	0-0.3': ASPHALT. $0.3-2.0$ ': Sand, F and silt and gravel; dusky brown (5YR 4/4), dry.	0 0	AS	00 00	
-2-	2-	N/A	/	2222	Same as above.	0 0		00 00 00 00 00 00 00 00 00 00 00 00 00	0 ppm
-4 <i>-</i>	4 —	N/A		3 3 2 4	Same as above, moist.	0 0	SW		0 ppm
-6 -	6 —	-05	/	7 8 7 9	Same as above, wet.	0.0000			0 ppm
-8-	8-	N/A	/	9 15 21 34	Sand, F-M and silt; trace gravel and C sand; moderate brown (5YR 4/4), wet.	o o			0 ppm
-10 <i>-</i>	10 —	N/A	/	14 15 12 12	Same as above.		SP		0 ppm
-12 <i>-</i>	12 —	N/A		15 14 14 19	Silt and clay; pale olive, wet.		ML		0 ppm
-14 -	14 —				End of boring at 14 feet.				
-16-	16-								
Chec	cked B	y: BEI	K_					Page 1 of	1



APPENDIX C

PREMIER LABORATORY CERTIFICATES OF ANALYSIS, FUSS & O'NEILL DATA VERIFICATION NARRATIVES AND CERTIFICATIONS, AND DATA VALIDATION COMPLETION WORKSHEETS



Modified Tier II Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E710I54
Date Samples Received at Laboratory:	10/26/2007
Date of Review:	12/18/2007

Four sediment samples, including a field duplicate, were and submitted to Premier Laboratory in Dayville, Connecticut for analysis of semi-volatile organic compounds (SVOCs) by EPA Method 8270C. Dedicated sampling equipment was used and volatile organic compounds (VOCs) were not constituents of concern; therefore, no equipment or trip blanks were indicated.

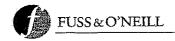
The relative percent differences (RPD) calculated for the primary and duplicate samples for several compounds were greater than the 50% limit established by the QAPP for non-aqueous samples.

Surrogate recoveries were acceptable for all samples. Low recovery of 2,4-dinitrophenol and high recovery of carbazole were reported for the lab control sample and/or lab control sample duplicate (LCS/LCSD). Reporting limits exceeded the Threshold Effects Concentrations (TECs) established by MacDonald, et al due to matrix interferences.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:

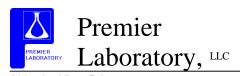
Lynne P. Matteso QA/QC Officer



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

		<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREMED	NTS:	,	_
Field measurement calibration records			N/
Groundwater field measurements (if applicable)			
Soil sampling field measurements (if applicable)			
Sediment sampling field measurements (if applicable)		₽	
Surface water sampling field measurements (if applicab	le)		□ 5\V
Low-flow sampling field measurements (if applicable)			Ala 🗆
Documentation of field activities		☑′_	
Sample numbering and labeling		<u> </u>	
Chain-of-Custody records		U	
Trip blanks			□ NA
Duplicate samples			□ N.
Equipment blanks			
Split samples (if any)			□ N/A
2. LABORATORY MEASUREMENTS:			
Trip blanks			□ \ A
Instrument blanks		⊠	
Laboratory control samples			
Duplicates samples			
Equipment blanks			□ ×
Matrix spike/matrix spike duplicates			□ v/\
Analysis type			
Chain-of-Custody records			
Surrogate recoveries			
Sample Project Narratives			
Split samples (if any)			
	TOTAL: _	13	
	PERCENT CO	ገለመ፣ ይሞይ.	100 %
	エニアヘニコット へん	وشلط بشارسا بالمالات	



61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

ANALYTICAL DATA REPORT

Report Number: E710I54 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

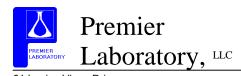
Received Date: 10/26/2007 Report Date: 11/5/2007

> Premier Laboratory, LLC Authorized Signature



Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

Page 1 of 18



61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

	MADEP MCP Analytical Method Report Certification Form								
Labo	ratory Name: Prei	mier Laboratory, LLC			Project #:	E710	54		
Proje	ct Location: Frank	klin, MA			MADEP RT	N ¹ :			
This	Form provides cert	ifications for the follow	ving data set:[list La	aboratory S	ample ID Num	ber(s)			
1, 2,	1, 2, 3, 4								
Samp	ole Matrices: 🗆 C	Groundwater ⊠ Soil/	Sediment 🗆 Drinl	king Water	☐ Other				
MCP	SW-846	8260B □	8151A □	8330	□ 6010	ов □	7		0A/1A □
Meth	ods Used	8270C ⊠	8081A □	VPH I		20 🗆		90)14M ² □
As spe	ecified in MADEP	8082 🗆	8021B □	EPH [[]	□ 7000 S	$S^3 \square$		7	7196A 🗆
	endium of	1 List Release Tracking No. 2 M - SW-846 Method 901	umber (RTN), if known	المماني ٨٠٠مناممان	o Cyanida (DAC)	Mathad			
	ical Methods.	3 S - SW-846 Methods 70				Metriou			
•		oonse to questions A				e Cert	ainty	" Si	tatus
Α		received by the labora				X \			No ¹
	1	the Chain-of-Custody							
В									No ¹
	included in this report followed, including the requirement to note and								
	discuss in a narrative QC data that did not meet appropriate performance								
	standards or guidelines?								
С									
	for "Presumptive	Certainty", as describe	ed in Section 2.0 (a),(b),(c) and	d (d) of the				
		t CAM VII A, "Quality	•	, , , , ,	` '				
		and Reporting of Ana		•					
D	VPH and EPH Me	ethods only: Was the	e VPH or EPH metl	hod run with	nout	`	Yes		No ¹
	significant modific	ations, as specified in	Section 11.3?						
	A response	to questions E and F	below is required	d for "Pres	umptive Certa	ainty"	statı	ıs	
Е	Were all QC perfo	ormance standards an	d recommendation	s for the		_ \ \	Yes	X	No ¹
	specified methods	s achieved?							
F	Were results for a	II analyte-list compou	nds/elements for th	e specified		X \	Yes		No ¹
	method(s) reporte	ed?							
	¹ All NO answ	ers must be addresse	d in an attached Er	nvironmenta	l Laboratory c	ase na	rrativ	/e.	
		der the pains and penalties for obtaining the information			nal				
inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.									
Sign	Signature: Position: Laboratory Director								
Print	ed Name: Rober	t Stevenson		Date: 11/5	5/2007				
II .									

Page 2 of 18

61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

> Report No: E710I54 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory, LLC received four samples from Fuss & O'Neill on 10/26/2007. The samples were analyzed from the following list of analyses:

Moisture, Percent

Semivolatiles by 8270C for GW/SW 8270C[3500]

Variances:

SDG:

2,4 - Dinitrophenol and Carbazole recoveries in the LCS/LCS duplicate were outside quality control limits. No action was necessary because the failures were less than 20% of all analytes, as specified in MCP Method 8270. These compounds were not detected in the sample.

Method:

None reported.

QA/QC:

Sample 3, 937071025-03, Semivolatiles by SW-846 8270C: One internal standard was below quality control limits for the sample due to matrix interference.

Sample 3, 937071025-03, Semivolatiles by SW-846 8270C: The detection limits are elevated for the sample due to matrix interference, as there was a large fuel pattern present.

Sample 4, 937071025-04, Semivolatiles by SW-846 8270C: One internal standard was below quality control limits for the sample due to matrix interference.

Sample 4, 937071025-04, Semivolatiles by SW-846 8270C: The detection limits are elevated for the sample due to matrix interference, as there was a large fuel pattern present.

Page 3 of 18

Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 1

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD Method: 8270C

QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-01

Matrix: Solid

Percent Moisture: 13.9

Sample Weight/Volume: 30.23 g

Dilution Factor: 5
Extract Volume: 1
Lab Data File: L20636.D

CAS No.	Parameter	Result	DL
103-33-3	Azobenzene	ND	960
83-32-9	Acenaphthene	ND	38
208-96-8	Acenaphthylene	77	38
62-53-3	Aniline	ND	1900
120-12-7	Anthracene	340	38
92-52-4	Biphenyl	ND	960
56-55-3	Benzo[a]anthracene	440	38
50-32-8	Benzo[a]pyrene	330	38
205-99-2	Benzo[b]fluoranthene	470	38
191-24-2	Benzo[g,h,i]perylene	ND	38
207-08-9	Benzo[k]fluoranthene	390	38
65-85-0	Benzoic acid	ND	4800
100-51-6	Benzyl alcohol	ND	1900
85-68-7	Benzyl butyl phthalate	ND	960
111-91-1	Bis(2-chloroethoxy)methane	ND	960
111-44-4	Bis(2-chloroethyl)ether	ND	960
108-60-1	Bis(2-chloroisopropyl)ether	ND	1900
117-81-7	Bis(2-ethylhexyl)phthalate	ND	960
101-55-3	4-Bromophenyl phenyl ether	ND	960
59-50-7	4-Chloro-3-methylphenol	ND	960
106-47-8	4-Chloroaniline	ND	1900
91-58-7	2-Chloronaphthalene	ND	960
95-57-8	2-Chlorophenol	ND	960
7005-72-3	4-Chlorophenyl phenyl ether	ND	960
218-01-9	Chrysene	480	38
53-70-3	Dibenz[a,h]anthracene	46	38
84-74-2	Di-n-butyl phthalate	ND	960
117-84-0	Di-n-octyl phthalate	ND	960
132-64-9	Dibenzofuran	ND	1900
95-50-1	1,2-Dichlorobenzene	ND	960
541-73-1	1,3-Dichlorobenzene	ND	960
106-46-7	1,4-Dichlorobenzene	ND	960
91-94-1	3,3-Dichlorobenzidine	ND	960
120-83-2	2,4-Dichlorophenol	ND	960
84-66-2	Diethyl phthalate	ND	960
131-11-3	Dimethyl phthalate	ND	960
105-67-9	2,4-Dimethylphenol	ND	960
51-28-5	2,4-Dinitrophenol	ND	960
121-14-2	2,4-Dinitrotoluene	ND	960
606-20-2	2,6-Dinitrotoluene	ND	960
206-44-0	Fluoranthene	890	38

6CSD9E50 Page 4 of 18

Laboratory: Premier Laboratory, LLC

PL Report No: E710I54
PL Sample No: 1 (continued)

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD

Method: 8270C QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-01

Matrix: Solid

Percent Moisture: 13.9

Sample Weight/Volume: 30.23 g

Dilution Factor: 5
Extract Volume: 1
Lab Data File: L20636.D

CAS No.	Parameter	Result	DL
86-73-7	Fluorene	ND	38
118-74-1	Hexachlorobenzene	ND	960
87-68-3	Hexachlorobutadiene	ND	960
77-47-4	Hexachlorocyclopentadiene	ND	960
67-72-1	Hexachloroethane	ND	960
193-39-5	Indeno[1,2,3-cd]pyrene	110	38
78-59-1	Isophorone	ND	960
534-52-1	2-Methyl-4,6-dinitrophenol	ND	960
91-57-6	2-Methylnaphthalene	ND	960
95-48-7	2-Methylphenol	ND	960
	3- & 4-Methylphenols	ND	960
91-20-3	Naphthalene	ND	38
88-74-4	2-Nitroaniline	ND	1900
99-09-2	3-Nitroaniline	ND	1900
100-01-6	4-Nitroaniline	ND	1900
98-95-3	Nitrobenzene	ND	960
88-75-5	2-Nitrophenol	ND	960
100-02-1	4-Nitrophenol	ND	960
621-64-7	N-Nitrosodi-n-propylamine	ND	960
62-75-9	N-Nitrosodimethylamine	ND	960
86-30-6	N-Nitrosodiphenylamine	ND	960
87-86-5	Pentachlorophenol	ND	960
85-01-8	Phenanthrene	ND	38
108-95-2	Phenol	ND	960
129-00-0	Pyrene	740	38
120-82-1	1,2,4-Trichlorobenzene	ND	960
95-95-4	2,4,5-Trichlorophenol	ND	960
88-06-2	2,4,6-Trichlorophenol	ND	960
Surrogate	Pacovary	Limite	

Surrogate	Recovery	Limits
2,4,6-Tribromophenol	64%	30%-130%
2-Fluorobiphenyl	49%	30%-130%
2-Fluorophenol	40%	30%-130%
4-Terphenyl-d14	79%	30%-130%
Nitrobenzene-d5	37%	30%-130%
Phenol-d6	42%	30%-130%

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 2

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD Method: 8270C

QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-02

Matrix: Solid

Percent Moisture: 17.6

Sample Weight/Volume: 30.59 g

Dilution Factor: 5 Extract Volume: 1 Lab Data File: L20637.D

103-33-3	CAS No.	Parameter	Result	DL
208-96-8 Acenaphthylene 160 40 62-53-3 Aniline ND 2000 120-12-7 Anthracene 280 40 92-52-4 Biphenyl ND 990 56-55-3 Benzo[alphyrene ND 40 50-32-8 Benzo[alpyrene 770 40 205-99-2 Benzo[billouranthene 690 40 191-24-2 Benzo[shilperylene ND 40 207-08-9 Benzo[killouranthene 1000 40 65-85-0 Benzole acid ND 5000 100-51-6 Benzyl alcohol ND 2000 85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chloroethoxylmethane ND 990 117-81-7 Benzyl butyl phthalate ND 990 117-81-7 Bis(2-chloroethy)ether ND 990 108-60-1 Bis(2-chloroethy)ether ND 990 108-60-1 Bis(2-chloroethy)ether ND 990	103-33-3	Azobenzene	ND	990
62-53-3	83-32-9	Acenaphthene	ND	40
120-12-7	208-96-8	Acenaphthylene	160	40
92-52-4 Biphenyl ND 990 56-55-3 Benzo[a]pyrene 770 40 205-99-2 Benzo[b]fluoranthene 690 40 205-99-2 Benzo[b]fluoranthene 690 40 191-24-2 Benzo[k]fluoranthene 1000 40 207-08-9 Benzo[k]fluoranthene 1000 40 65-85-0 Benzyl alcohol ND 5000 100-51-6 Benzyl alcohol ND 2000 85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chlorocthyz)ether ND 990 118-60-1 Bis(2-chlorocthyz)ether ND 990 118-71-7 Bis(2-chlorospropybether ND 2000 117-81-7 Bis(2-chlorospropybether ND 2000 117-81-7 Bis(2-chlorospropybether ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 105-47-8 4-Chloro-3-methylphenol ND 990 95-57-8 2-Chloropheny	62-53-3	Aniline	ND	2000
56-55-3 Benzo[a]anthracene ND 40 50-32-8 Benzo[a]pyrene 770 40 205-99-2 Benzo[b]fluoranthene 690 40 191-24-2 Benzo[g,h,i]perylene ND 40 207-08-9 Benzo[g,h,i]perylene ND 40 65-85-0 Benzola caid ND 5000 100-51-6 Benzyl alcohol ND 2000 85-68-7 Benzyl alcohol ND 990 111-91-1 Bis(2-chloroethylene ND 990 111-44-4 Bis(2-chloroethyl)ether ND 990 108-60-1 Bis(2-chloroethyl)ether ND 990 108-60-1 Bis(2-chloroisopropylether ND 990 107-81-7 Bis(2-chloroisopropylether ND 990 108-60-1 Bis(2-chloroisopropylether ND 990 108-7-8 4-Bromophenyl phenyl ether ND 990 108-5-3 4-Bromophenyl phenyl ether ND 990 91-58-7 2-Chloroalili	120-12-7	Anthracene	280	40
56-55-3 Benzo[a]anthracene ND 40 50-32-8 Benzo[a]pyrene 770 40 205-99-2 Benzo[b]fluoranthene 690 40 191-24-2 Benzo[g,h,i]perylene ND 40 207-08-9 Benzo[g,h,i]perylene ND 40 65-85-0 Benzo[a caid ND 5000 100-51-6 Benzyl alcohol ND 900 85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chlorocethoxy)methane ND 990 111-44-4 Bis(2-chlorocethy)bether ND 990 117-81-7 Bis(2-chlorocethy)bether ND 990 108-60-1 Bis(2-chlorocethy)bether ND 990 108-57-3 4-Bromophenyl phenyl ether ND 990 104-78-8 4-Chloro-3-methylphenol ND 990 91-58-7 2-Chlorophenyl phenyl ether ND 990 95-57-8 2-Chlorophenol ND 990 95-57-8 2-Chloropheny	92-52-4	Biphenyl	ND	990
Description	56-55-3	Benzo[a]anthracene	ND	40
Description	50-32-8	Benzo[a]pyrene	770	40
Benzo[k]fluoranthene 1000 40	205-99-2		690	40
207-08-9 Benzoic acid ND 5000 65-85-0 Benzoic acid ND 5000 100-51-6 Benzyl alcohol ND 2000 85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chloroethoxy)methane ND 990 111-44-4 Bis(2-chloroethy)lether ND 990 108-60-1 Bis(2-chloroethy)lether ND 990 117-81-7 Bis(2-chlorospopt)lether ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 106-47-8 4-Chloro-3-methylphenol ND 990 91-58-7 2-Chloroaphthalene ND 990 95-57-8 2-Chlorophenyl phenyl ether ND 990 95-57-8 2-Chlorophenyl phenyl ether ND 990 18-01-9 Chrysene 1100 40 18-4-42 Di-n-butyl phthalate ND 990 117-84-0 Di-n-	191-24-2	Benzo[g,h,i]perylene	ND	40
100-51-6 Benzyl alcohol ND 2000 85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chloroethoxy)methane ND 990 111-44-4 Bis(2-chloroethy)lether ND 990 108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 106-47-8 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chloroaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 95-57-8 2-Chlorophenol ND 990 128-01-9 Chrysene 1100 40 40 40 40 40 40	207-08-9	Benzo[k]fluoranthene	1000	40
85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chloroethoxy)methane ND 990 111-44-4 Bis(2-chloroethy)lether ND 990 108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 990 91-58-7 2-Chlorophenol ND 990 95-57-8 2-Chlorophenol ND 990 218-01-9 Chrysene 1100 40 218-01-9 Chrysene 1100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene	65-85-0	Benzoic acid	ND	5000
85-68-7 Benzyl butyl phthalate ND 990 111-91-1 Bis(2-chloroethoxy)methane ND 990 111-44-4 Bis(2-chloroethoxy)methane ND 990 108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chlorophenol ND 990 95-57-8 2-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 218-01-9 Chrysene 1100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 55-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dich	100-51-6	Benzyl alcohol	ND	2000
111-91-1 Bis(2-chloroethoxy)methane ND 990 111-44-4 Bis(2-chloroethyl)ether ND 990 108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroanline ND 990 91-58-7 2-Chloronaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-butyl phthalate ND 990 132-64-9 Dibenzofuran ND 990 541-73-1 1,3-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 120-83-2 2,4-Dichlor	85-68-7		ND	990
111-44-4 Bis(2-chloroethyl)ether ND 990 108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 990 91-58-7 2-Chlorophenol ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 990 55-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene	111-91-1		ND	990
108-60-1 Bis(2-chloroisopropyl)ether ND 2000 117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaliline ND 2000 91-58-7 2-Chlorophenol ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-cotyl phthalate ND 990 132-64-9 Dibenzofuran ND 990 55-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorobenzidine <td>111-44-4</td> <td></td> <td>ND</td> <td>990</td>	111-44-4		ND	990
117-81-7 Bis(2-ethylhexyl)phthalate ND 990 101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chloroaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzidine ND 990 106-46-7 1,4-Dichlorobenzidine ND 990 12-83-2 2,4-Dichlorobenzidine	108-60-1		ND	2000
101-55-3 4-Bromophenyl phenyl ether ND 990 59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chlorophenol ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-cytl phthalate ND 990 122-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 12-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND			ND	
59-50-7 4-Chloro-3-methylphenol ND 990 106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chloronaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-butyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND <td></td> <td></td> <td></td> <td></td>				
106-47-8 4-Chloroaniline ND 2000 91-58-7 2-Chloronaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzene ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 13-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dimitrophenol ND 990		1 1 1		990
91-58-7 2-Chloronaphthalene ND 990 95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				2000
95-57-8 2-Chlorophenol ND 990 7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 990 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				
7005-72-3 4-Chlorophenyl phenyl ether ND 990 218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				990
218-01-9 Chrysene 1100 40 53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				990
53-70-3 Dibenz[a,h]anthracene 100 40 84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				40
84-74-2 Di-n-butyl phthalate ND 990 117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990				
117-84-0 Di-n-octyl phthalate ND 990 132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	84-74-2		ND	990
132-64-9 Dibenzofuran ND 2000 95-50-1 1,2-Dichlorobenzene ND 990 541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	117-84-0		ND	990
541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	132-64-9		ND	2000
541-73-1 1,3-Dichlorobenzene ND 990 106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	95-50-1	1,2-Dichlorobenzene	ND	990
106-46-7 1,4-Dichlorobenzene ND 990 91-94-1 3,3-Dichlorobenzidine ND 990 120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990			ND	990
120-83-2 2,4-Dichlorophenol ND 990 84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	106-46-7	1,4-Dichlorobenzene	ND	990
84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	91-94-1	3,3-Dichlorobenzidine	ND	990
84-66-2 Diethyl phthalate ND 990 131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	120-83-2	2,4-Dichlorophenol	ND	990
131-11-3 Dimethyl phthalate ND 990 105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	84-66-2		ND	990
105-67-9 2,4-Dimethylphenol ND 990 51-28-5 2,4-Dinitrophenol ND 990	131-11-3		ND	990
51-28-5 2,4-Dinitrophenol ND 990	105-67-9	* *	ND	990
	51-28-5		ND	990
121-14-2 2,4-Dinitrotoluene ND 990				990
606-20-2 2,6-Dinitrotoluene ND 990				990
206-44-0 Fluoranthene 2500 40		•		

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 2 (continued)

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD

Method: 8270C QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-02

Matrix: Solid

Percent Moisture: 17.6

Sample Weight/Volume: 30.59 g

Dilution Factor: 5 Extract Volume: 1 Lab Data File: L20637.D

CAS No.	Parameter	Result	DL
86-73-7	Fluorene	81	40
118-74-1	Hexachlorobenzene	ND	990
87-68-3	Hexachlorobutadiene	ND	990
77-47-4	Hexachlorocyclopentadiene	ND	990
67-72-1	Hexachloroethane	ND	990
193-39-5	Indeno[1,2,3-cd]pyrene	240	40
78-59-1	Isophorone	ND	990
534-52-1	2-Methyl-4,6-dinitrophenol	ND	990
91-57-6	2-Methylnaphthalene	ND	990
95-48-7	2-Methylphenol	ND	990
	3- & 4-Methylphenols	ND	990
91-20-3	Naphthalene	ND	40
88-74-4	2-Nitroaniline	ND	2000
99-09-2	3-Nitroaniline	ND	2000
100-01-6	4-Nitroaniline	ND	2000
98-95-3	Nitrobenzene	ND	990
88-75-5	2-Nitrophenol	ND	990
100-02-1	4-Nitrophenol	ND	990
621-64-7	N-Nitrosodi-n-propylamine	ND	990
62-75-9	N-Nitrosodimethylamine	ND	990
86-30-6	N-Nitrosodiphenylamine	ND	990
87-86-5	Pentachlorophenol	ND	990
85-01-8	Phenanthrene	1300	40
108-95-2	Phenol	ND	990
129-00-0	Pyrene	2100	40
120-82-1	1,2,4-Trichlorobenzene	ND	990
95-95-4	2,4,5-Trichlorophenol	ND	990
88-06-2	2,4,6-Trichlorophenol	ND	990
Surrogate	Recovery	Limits	

Surrogate	Recovery	Limits
2,4,6-Tribromophenol	68%	30%-130%
2-Fluorobiphenyl	61%	30%-130%
2-Fluorophenol	50%	30%-130%
4-Terphenyl-d14	94%	30%-130%
Nitrobenzene-d5	50%	30%-130%
Phenol-d6	52%	30%-130%

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 3

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD Method: 8270C

QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-03

Matrix: Solid

Percent Moisture: 27.8

Sample Weight/Volume: 30.29 g

Dilution Factor: 5 Extract Volume: 1 Lab Data File: L20638.D

CAS No.	Parameter	Result	DL
103-33-3	Azobenzene	ND	1100
83-32-9	Acenaphthene	ND	46
208-96-8	Acenaphthylene	170	46
62-53-3	Aniline	ND	2300
120-12-7	Anthracene	98	46
92-52-4	Biphenyl	ND	1100
56-55-3	Benzo[a]anthracene	920	46
50-32-8	Benzo[a]pyrene	860	46
205-99-2	Benzo[b]fluoranthene	1200	46
191-24-2	Benzo[g,h,i]perylene	ND	46
207-08-9	Benzo[k]fluoranthene	990	46
65-85-0	Benzoic acid	ND	5700
100-51-6	Benzyl alcohol	ND	2300
85-68-7	Benzyl butyl phthalate	ND	1100
111-91-1	Bis(2-chloroethoxy)methane	ND	1100
111-44-4	Bis(2-chloroethyl)ether	ND	1100
108-60-1	Bis(2-chloroisopropyl)ether	ND	2300
117-81-7	Bis(2-ethylhexyl)phthalate	ND	1100
101-55-3	4-Bromophenyl phenyl ether	ND	1100
59-50-7	4-Chloro-3-methylphenol	ND	1100
106-47-8	4-Chloroaniline	ND	2300
91-58-7	2-Chloronaphthalene	ND	1100
95-57-8	2-Chlorophenol	ND	1100
7005-72-3	4-Chlorophenyl phenyl ether	ND	1100
218-01-9	Chrysene	1200	46
53-70-3	Dibenz[a,h]anthracene	160	46
84-74-2	Di-n-butyl phthalate	ND	1100
117-84-0	Di-n-octyl phthalate	ND	1100
132-64-9	Dibenzofuran	ND	2300
95-50-1	1,2-Dichlorobenzene	ND	1100
541-73-1	1,3-Dichlorobenzene	ND	1100
106-46-7	1,4-Dichlorobenzene	ND	1100
91-94-1	3,3-Dichlorobenzidine	ND	1100
120-83-2	2,4-Dichlorophenol	ND	1100
84-66-2	Diethyl phthalate	ND	1100
131-11-3	Dimethyl phthalate	ND	1100
105-67-9	2,4-Dimethylphenol	ND	1100
51-28-5	2,4-Dinitrophenol	ND	1100
121-14-2	2,4-Dinitrotoluene	ND	1100
606-20-2	2,6-Dinitrotoluene	ND	1100
206-44-0	Fluoranthene	2000	46

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 3 (continued)

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD

Method: 8270C QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-03

Matrix: Solid

Percent Moisture: 27.8

Sample Weight/Volume: 30.29 g

Dilution Factor: 5
Extract Volume: 1

Lab Data File: L20638.D

CAS No.	Parameter	Result	DL
86-73-7	Fluorene	ND	46
118-74-1	Hexachlorobenzene	ND	1100
87-68-3	Hexachlorobutadiene	ND	1100
77-47-4	Hexachlorocyclopentadiene	ND	1100
67-72-1	Hexachloroethane	ND	1100
193-39-5	Indeno[1,2,3-cd]pyrene	310	46
78-59-1	Isophorone	ND	1100
534-52-1	2-Methyl-4,6-dinitrophenol	ND	1100
91-57-6	2-Methylnaphthalene	ND	1100
95-48-7	2-Methylphenol	ND	1100
	3- & 4-Methylphenols	ND	1100
91-20-3	Naphthalene	ND	46
88-74-4	2-Nitroaniline	ND	2300
99-09-2	3-Nitroaniline	ND	2300
100-01-6	4-Nitroaniline	ND	2300
98-95-3	Nitrobenzene	ND	1100
88-75-5	2-Nitrophenol	ND	1100
100-02-1	4-Nitrophenol	ND	1100
621-64-7	N-Nitrosodi-n-propylamine	ND	1100
62-75-9	N-Nitrosodimethylamine	ND	1100
86-30-6	N-Nitrosodiphenylamine	ND	1100
87-86-5	Pentachlorophenol	ND	1100
85-01-8	Phenanthrene	ND	46
108-95-2	Phenol	ND	1100
129-00-0	Pyrene	2000	46
120-82-1	1,2,4-Trichlorobenzene	ND	1100
95-95-4	2,4,5-Trichlorophenol	ND	1100
88-06-2	2,4,6-Trichlorophenol	ND	1100
Surrogate	Recovery	Limits	

Surrogate	Recovery	Limits	
2,4,6-Tribromophenol	65%	30%-130%	
2-Fluorobiphenyl	76%	30%-130%	
2-Fluorophenol	66%	30%-130%	
4-Terphenyl-d14	99%	30%-130%	
Nitrobenzene-d5	58%	30%-130%	
Phenol-d6	68%	30%-130%	

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 4

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD Method: 8270C

QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-04

Matrix: Solid

Percent Moisture: 17.0

Sample Weight/Volume: 30.13 g

Dilution Factor: 5
Extract Volume: 1
Lab Data Files J 200

Lab Data File: L20639.D

CAS No.	Parameter	Result	DL
103-33-3	Azobenzene	ND	1000
83-32-9	Acenaphthene	ND	40
208-96-8	Acenaphthylene	76	40
62-53-3	Aniline	ND	2000
120-12-7	Anthracene	58	40
92-52-4	Biphenyl	ND	1000
56-55-3	Benzo[a]anthracene	400	40
50-32-8	Benzo[a]pyrene	350	40
205-99-2	Benzo[b]fluoranthene	520	40
191-24-2	Benzo[g,h,i]perylene	ND	40
207-08-9	Benzo[k]fluoranthene	450	40
65-85-0	Benzoic acid	ND	5000
100-51-6	Benzyl alcohol	ND	2000
85-68-7	Benzyl butyl phthalate	ND	1000
111-91-1	Bis(2-chloroethoxy)methane	ND	1000
111-44-4	Bis(2-chloroethyl)ether	ND	1000
108-60-1	Bis(2-chloroisopropyl)ether	ND	2000
117-81-7	Bis(2-ethylhexyl)phthalate	ND	1000
101-55-3	4-Bromophenyl phenyl ether	ND	1000
59-50-7	4-Chloro-3-methylphenol	ND	1000
106-47-8	4-Chloroaniline	ND	2000
91-58-7	2-Chloronaphthalene	ND	1000
95-57-8	2-Chlorophenol	ND	1000
7005-72-3	4-Chlorophenyl phenyl ether	ND	1000
218-01-9	Chrysene	500	40
53-70-3	Dibenz[a,h]anthracene	ND	40
84-74-2	Di-n-butyl phthalate	ND	1000
117-84-0	Di-n-octyl phthalate	ND	1000
132-64-9	Dibenzofuran	ND	2000
95-50-1	1,2-Dichlorobenzene	ND	1000
541-73-1	1,3-Dichlorobenzene	ND	1000
106-46-7	1,4-Dichlorobenzene	ND	1000
91-94-1	3,3-Dichlorobenzidine	ND	1000
120-83-2	2,4-Dichlorophenol	ND	1000
84-66-2	Diethyl phthalate	ND	1000
131-11-3	Dimethyl phthalate	ND	1000
105-67-9	2,4-Dimethylphenol	ND	1000
51-28-5	2,4-Dinitrophenol	ND	1000
121-14-2	2,4-Dinitrotoluene	ND	1000
606-20-2	2,6-Dinitrotoluene	ND	1000
206-44-0	Fluoranthene	810	40

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Laboratory: Premier Laboratory, LLC

PL Report No: E710I54 PL Sample No: 4 (continued)

Date Collected: 10/25/2007 Date Received: 10/26/2007 Date Extracted: 10/31/07 By: KT

Date Analyzed: 11/01/07 By: JD

Method: 8270C QC Batch#: 57560 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 937071025-04

Matrix: Solid

Percent Moisture: 17.0

Sample Weight/Volume: 30.13 g

Dilution Factor: 5 Extract Volume: 1 Lab Data File: L20639.D

CAS No.	Parameter	Result	DL
86-73-7	Fluorene	ND	40
118-74-1	Hexachlorobenzene	ND	1000
87-68-3	Hexachlorobutadiene	ND	1000
77-47-4	Hexachlorocyclopentadiene	ND	1000
67-72-1	Hexachloroethane	ND	1000
193-39-5	Indeno[1,2,3-cd]pyrene	130	40
78-59-1	Isophorone	ND	1000
534-52-1	2-Methyl-4,6-dinitrophenol	ND	1000
91-57-6	2-Methylnaphthalene	ND	1000
95-48-7	2-Methylphenol	ND	1000
	3- & 4-Methylphenols	ND	1000
91-20-3	Naphthalene	ND	40
88-74-4	2-Nitroaniline	ND	2000
99-09-2	3-Nitroaniline	ND	2000
100-01-6	4-Nitroaniline	ND	2000
98-95-3	Nitrobenzene	ND	1000
88-75-5	2-Nitrophenol	ND	1000
100-02-1	4-Nitrophenol	ND	1000
621-64-7	N-Nitrosodi-n-propylamine	ND	1000
62-75-9	N-Nitrosodimethylamine	ND	1000
86-30-6	N-Nitrosodiphenylamine	ND	1000
87-86-5	Pentachlorophenol	ND	1000
85-01-8	Phenanthrene	ND	40
108-95-2	Phenol	ND	1000
129-00-0	Pyrene	950	40
120-82-1	1,2,4-Trichlorobenzene	ND	1000
95-95-4	2,4,5-Trichlorophenol	ND	1000
88-06-2	2,4,6-Trichlorophenol	ND	1000
Surrogate	Recovery	Limits	

Surrogate	Recovery	Limits	
2,4,6-Tribromophenol	68%	30%-130%	
2-Fluorobiphenyl	75%	30%-130%	
2-Fluorophenol	66%	30%-130%	
4-Terphenyl-d14	106%	30%-130%	
Nitrobenzene-d5	63%	30%-130%	
Phenol-d6	67%	30%-130%	

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FORM 3 Soil 8270C Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20632.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	()	()	Rec#	Rec
1,2,4-Trichlorobenzene	666.7	444.3	67	30-130
1,2-Dichlorobenzene	666.7	384.9	58	30-130
1,3-Dichlorobenzene	666.7	357.6	54	30-130
1,4-Dichlorobenzene	666.7	349.2	52	30-130
2,4,5-Trichlorophenol	666.7	490.3	74	30-130
2,4,6-Trichlorophenol	666.7	499.7	75	30-130
2,4-Dichlorophenol	666.7	467.2	70	30-130
2,4-Dimethylphenol	666.7	375.9	56	30-130
2,4-Dinitrophenol	666.7	116.3	17*	30-130
2,4-Dinitrotoluene	666.7	642.3	96	30-130
2,6-Dinitrotoluene	666.7	616.5	92	30-130
2-Chloronaphthalene	666.7	523.4	78	30-130
2-Chlorophenol	666.7	443.5	66	30-130
2-Methylnaphthalene	666.7	487.9	73	30-130
2-Methylphenol	666.7	461.3	69	30-130
2-Nitroaniline	666.7	573.8	86	30-130
2-Nitrophenol	666.7	392.0	59	30-130
3- & 4-Methylphenols	666.7	466.7	70	30-130
3-Nitroaniline	666.7	660.7	99	30-130
4-Chloroaniline	666.7	481.6	72	30-130
4-Chlorophenyl phenyl	666.7	550.2	82	30-130
4-Nitroaniline	666.7	657.8	99	30-130
4-Nitrophenol	666.7	716.4	107	30-130
Acenaphthene	666.7	537.9	81	30-130
Acenaphthylene	666.7	556.6	83	30-130
Aniline	666.7	422.0	63	30-130
Anthracene	666.7	706.8	106	30-130
Azobenzene	666.7	675.8	101	30-130

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil 8270C Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20632.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	()	()	Rec#	Rec
Benzo[a]pyrene	666.7	635.5	95	30-130
Benzo[b]fluoranthene	666.7	669.0	100	30-130
<pre>Benzo[g,h,i]perylene</pre>	666.7	666.0	100	30-130
Benzo[k]fluoranthene	666.7	638.3	96	30-130
Benzyl alcohol	666.7	455.9	68	30-130
Benzyl butyl phthalate	666.7	742.4	111	30-130
Bis(2-chloroethoxy)me	666.7	489.2	73	30-130
Bis(2-chloroethyl)ether	666.7	461.1	69	30-130
Bis(2-chloroisopropyl	666.7	394.6	59	30-130
Bis(2-ethylhexyl)phth	666.7	749.2	112	30-130
Carbazole	666.7	867.2	130	30-130
Chrysene	666.7	704.8	106	30-130
Di-n-butyl phthalate	666.7	698.5	105	30-130
Di-n-octyl phthalate	666.7	656.8	98	30-130
Dibenzofuran	666.7	569.2	85	30-130
Dibenz[a,h]anthracene	666.7	611.0	92	30-130
Diethyl phthalate	666.7	637.3	96	30-130
Dimethyl phthalate	666.7	574.4	86	30-130
Fluoranthene	666.7	772.5	116	30-130
Fluorene	666.7	601.5	90	30-130
Hexachlorobenzene	666.7	652.6	98	30-130
Hexachlorobutadiene	666.7	458.7	69	30-130
Hexachlorocyclopentad	666.7	551.7	83	30-130
Hexachloroethane	666.7	364.9	55	30-130
<pre>Indeno[1,2,3-cd]pyrene</pre>	666.7	614.0	92	30-130
Isophorone	666.7	497.3	74	30-130
N-Nitrosodi-n-propyla	666.7	460.5	69	30-130
N-Nitrosodimethylamine	666.7	340.5	51	30-130

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil 8270C Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20632.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	()	()	Rec#	Rec
N-Nitrosodiphenylamine	666.7	758.6	114	30-130
Naphthalene	666.7	486.8	73	30-130
Nitrobenzene	666.7	459.6	69	30-130
Pentachlorophenol	666.7	544.8	82	30-130
Phenanthrene	666.7	747.1	112	30-130
Phenol	666.7	468.6	70	30-130
Pyrene	666.7	706.6	106	30-130

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil 8270C Lab Control Sample Duplicate

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20633.D

	Spike	Sample				QC
	Added	Concentration	%		Li	mits
Compound	()	()	Rec#	RPD#	RPD	Rec
1,2,4-Trichlorobenzene	666.7	467.4	70	4.38	30	30-130
1,2-Dichlorobenzene	666.7	413.6	62	6.67	30	30-130
1,3-Dichlorobenzene	666.7	392.5	59	8.85	30	30-130
1,4-Dichlorobenzene	666.7	395.2	59	12.6	30	30-130
2,4,5-Trichlorophenol	666.7	574.8	86	15.0	30	30-130
2,4,6-Trichlorophenol	666.7	550.0	82	8.92	30	30-130
2,4-Dichlorophenol	666.7	465.9	70	0	30	30-130
2,4-Dimethylphenol	666.7	361.2	54	3.64	30	30-130
2,4-Dinitrophenol	666.7	114.2	17*	0	30	30-130
2,4-Dinitrotoluene	666.7	686.0	103	7.04	30	30-130
2,6-Dinitrotoluene	666.7	647.6	97	5.29	30	30-130
2-Chloronaphthalene	666.7	551.5	83	6.21	30	30-130
2-Chlorophenol	666.7	460.7	69	4.44	30	30-130
2-Methylnaphthalene	666.7	518.4	78	6.62	30	30-130
2-Methylphenol	666.7	494.9	74	6.99	30	30-130
2-Nitroaniline	666.7	616.5	92	6.74	30	30-130
2-Nitrophenol	666.7	424.7	64	8.13	30	30-130
3- & 4-Methylphenols	666.7	510.8	77	9.52	30	30-130
3-Nitroaniline	666.7	673.3	101	2.00	30	30-130
4-Chloroaniline	666.7	492.2	74	2.74	30	30-130
4-Chlorophenyl phenyl	666.7	561.0	84	2.41	30	30-130
4-Nitroaniline	666.7	694.6	104	4.93	30	30-130
4-Nitrophenol	666.7	584.0	88	19.5	30	30-130
Acenaphthene	666.7	568.2	85	4.82	30	30-130
Acenaphthylene	666.7	595.4	89	6.98	30	30-130
Aniline	666.7	471.2		11.9	30	30-130
Anthracene	666.7	754.5	113	6.39	30	30-130
Azobenzene	666.7	733.8	110	8.53	30	30-130

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil 8270C Lab Control Sample Duplicate

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20633.D (continued)

	Spike	Sample				QC
	Added	Concentration	%		Li	mits
Compound	()	()	Rec#	RPD#	RPD	Rec
Benzo[a]pyrene	666.7	687.3		8.08	30	30-130
Benzo[b]fluoranthene	666.7	699.8	105	4.88	30	30-130
Benzo[g,h,i]perylene	666.7	726.5	109	8.61	30	30-130
Benzo[k]fluoranthene	666.7	681.7	102	6.06	30	30-130
Benzyl alcohol	666.7	491.5	74	8.45	30	30-130
Benzyl butyl phthalate	666.7	787.9	118	6.11	30	30-130
Bis(2-chloroethoxy)me	666.7	502.0	75	2.70	30	30-130
Bis(2-chloroethyl)ether	666.7	452.8	68	1.46	30	30-130
Bis(2-chloroisopropyl	666.7	428.7	64	8.13	30	30-130
Bis(2-ethylhexyl)phth	666.7	801.6	120	6.90	30	30-130
Carbazole	666.7	921.2	138*	5.97	30	30-130
Chrysene	666.7	762.8	114	7.27	30	30-130
Di-n-butyl phthalate	666.7	735.6	110	4.65	30	30-130
Di-n-octyl phthalate	666.7	688.4	103	4.98	30	30-130
Dibenzofuran	666.7	617.6	93	8.99	30	30-130
Dibenz[a,h]anthracene	666.7	644.4	97	5.29	30	30-130
Diethyl phthalate	666.7	668.1	100	4.08	30	30-130
Dimethyl phthalate	666.7	612.8	92	6.74	30	30-130
Fluoranthene	666.7	811.7	122	5.04	30	30-130
Fluorene	666.7	630.0	94	4.35	30	30-130
Hexachlorobenzene	666.7	691.4	104	5.94	30	30-130
Hexachlorobutadiene	666.7	486.2	73	5.63	30	30-130
Hexachlorocyclopentad	666.7	563.0	84	1.20	30	30-130
Hexachloroethane	666.7	382.2	57	3.57	30	30-130
<pre>Indeno[1,2,3-cd]pyrene</pre>	666.7	646.4	97	5.29	30	30-130
Isophorone	666.7	531.3		7.79	30	30-130
N-Nitrosodi-n-propyla	666.7	492.4		6.99	30	30-130
N-Nitrosodimethylamine	666.7	366.3	55	7.55	30	30-130

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil 8270C Lab Control Sample Duplicate

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/01/07

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1031S-1 Location: Franklin, MA

Lab File ID: L20633.D (continued)

	Spike	Sample				QC
	Added	Concentration	%		Limits	
Compound	()	()	Rec#	RPD#	RPD	Rec
N-Nitrosodiphenylamine	666.7	814.0	122	6.78	30	30-130
Naphthalene	666.7	511.9	77	5.33	30	30-130
Nitrobenzene	666.7	470.9	71	2.86	30	30-130
Pentachlorophenol	666.7	615.2	92	11.5	30	30-130
Phenanthrene	666.7	800.8	120	6.90	30	30-130
Phenol	666.7	461.2	69	1.44	30	30-130
Pyrene	666.7	759.8	114	7.27	30	30-130

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 4 8270C Method Blank Summary

Project No.: E710I54 Project: 20050458.B10/Nu-Style Phase

Lab File ID: L20634.D Lab Sample ID: S1031BS-1

Matrix: Soil Date Analyzed: 11/01/07

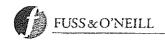
Instrument ID: MS10 Date Extracted:

Time Analyzed: 1406

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E710I54-1	937071025-01	L20636.D	11/01/2007
2	E710I54-2	937071025-02	L20637.D	11/01/2007
3	E710I54-3	937071025-03	L20638.D	11/01/2007
4	E710I54-4	937071025-04	L20639.D	11/01/2007
5	LCS1031S-1	LCS1031S-1	L20632.D	11/01/2007
6	LCSD1031s-1	LCSD1031S-1	L20633.D	11/01/2007
7				
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Page 18 of 18



INITIAL DATE: OCTOBER 2006 REVISION DATE: OCTOBER 2006 REVISION: 10

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

			YES	NO COMMENTS
1. 2. 3.	SDG Project Narratives Traffic Report Volatiles Data	B .		
	a. Sample Data Target Compound List (TCL) Results			MA
	Reconstructed total ion chromatograms (RIC) for each Sample For each sample:			N IB
	Raw spectra and background-subtracted mass spectra of target compounds identified Mass spectra of all reported TICs with three best library			NA
	matches Percent solids calculations	X		NIA
1	b. Standards Data (all instruments) Initial Calibration Data RICs and Quan Reports for all Standards Continuing Calibration RICs and Quan Reports for all Standards Internal Standard Area Summary			NIA NIA NIA NIA NIA
C	:. Raw QC Data Blank Data Matrix Spike Data Matrix Spike Duplicate Data			NIA NIA
S a.	emivolatiles Data . QC Summary Surrogate Percent Recovery Summary MS/MSD Summary		6 	compounds out in KS ess than 20%, meets MCP requirements VIA
	Method Blank Summary Tuning and Mass Calibration	生		

4.



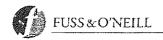
INITIAL DATE: OCTOBER 2006 REVISION DATE: OCTOBER 2006 REVISION: 1.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

YES NO COMMENTS

ъ.	Sample Data			
	TCL Results			
	Tentatively Identified Compounds			NIA
	Reconstructed total ion chromatograms (RIC) for each			
	Sample] .	
	For each sample:			
	Raw spectra and background-subtracted mass			
	spectra of TCL compounds] _	
	Mass spectra of TICs with 3 best library matches] _	NIA
	GPC chromatograms (if GPC performed)			NIN
c.				
	Initial Calibration Data	X C		
	RICs and Quan Reports for all Standards		_	
	Continuing Calibration	B	_	
	RICs and Quan Reports for all Standards			
	Internal Standard Areas Summary		_	
	Internal Standard Areas Summary] _	
	Raw QC Data			
	Decatluorotripbenylphosphine (DFTPP)] _	·····
	Blank Data		_	
	Matrix Spike Data		_	1114
	Matrix Spike Duplicate Data		_	WIU
	rellaneous Data			
	Original preparation and analysis forms or copies of preparation			
	and analysis log book pages		_	
I	nternal sample & sample extract transfer chain-of custody			
ľ	records			
3	Screening Records			NIA
-	All instrument output, including strip charts from screening			
	crivities (describe or list)			



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

				<u>YES</u>	<u>NO</u>	COMMENTS	
6.	S	stody Records Jample Log-in Sheet (Lab & DC1) Liscellaneous Shipping/Receiving Recor	ds (describe or list)				
7.	Internal Lab	Sample Transfer Records and Trackir Sheets (describe or list)	ng		Ο		
8.	Other Record	ds (describe or list)			O		
9.	Comments:						
	**	See laboratory Quality Assurance Pl	an for limits.				
	Completed (Lab)	by: Ly Mile Gr (Signature) (P	egon lante stated Name/Title)	Orfa Mani	nics ager	11 /5 /0 Date	7-
		the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the above information is true and accurate the ac					
	Certified by (Lab)	Mulial Attoristan (P) (Signature)	obert Stevencon cinted Name (Title)	1/60	b Pire	dor 11-5-0 Date	2



146 H	artford	Road,	Manchester,	CT	0604
 ~ ~ ~					

☐ 56 Quarry Road, Trumbull, CT 06611

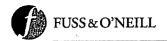
1419 Richland Street, Columbia, SC 29201

78 Interstate Drive, West Springfield, MA 01089
610 Lynndale Court, Suite E, Greenville, NC 27858

☐ 24 Madison Avenue Extension, Albany, NY 12203

275 Promenade Street, Suite 350, Providence, RI 02 08 U 80 Washington Street, Suite 301, Poughkeepsie, NY 12601

CHAIN-OF-CUSTODY RECORD	14715	Thrnaron 1 Day* U 3 Days*	Char (days)
PROJECT NAME PROJECT LOCATION Nu-Style Phase II Franklin, MA REPORT TO: David Foss, dfoss@fands.com Invoice To:	PROJECT NUMBER 2005 0 458 Analysis Request	310	LABORATORY Premier Containers
P.O. No.: \$4/2006 0458 BlO Sampler's Signature: Date: 10/25/07 Source Codes: W=Potable Water S=Soil W=Waste S=Surface Water T=Treatment Facility B=Sediment A=Air		/ //	
Item No. Transfer Check No. Sample Number Source Code Sampled Date Sampled Time Sampled 1 1 2 3 4 10/25/07 1430 10/25/07 1430 1435 1435 1440 1440 1440 1445 <td>/</td> <td> </td> <td>Comments</td>	/		Comments
Transfer Number Relinquished By Accepted By Date 1	MCP Dota Enhances Additional Comments:	uirements: ment Project - see offac 41.4 etc attached QM	



Modified Tier II Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E711143
Date Samples Received at Laboratory:	11/2/2007
Date of Review:	12/18/2007

Four soil samples, including a field duplicate, were collected from four soil borings over two days. Samples were submitted to Premier Laboratory in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant-13 metals plus barium by EPA Methods 6010B and 7471, and/or petroleum hydrocarbons by Massachusetts Extractable and Volatile Petroleum Hydrocarbon (EPH/VPH) methods of analysis. Analyses for individual samples were determined based upon constituents of concern in the area where the individual monitoring wells are located.

Dedicated sampling equipment was used; therefore, no equipment blank was indicated. An aqueous trip blank was also submitted. No VOCs were reported in the trip blank.

Results of primary and duplicate sample pairs were generally similar. The relative percent difference (RPD) calculated for several metals was greater than the 50% limit established by the QAPP for soil samples. Several RPD values were outside laboratory quality control limits in both initial total and fractionated extractions.

Surrogate recoveries were acceptable for all applicable analyses. With the exception of hexachlorobutadiene, reporting limits were low enough to compare to MADEP GW-2 and GW-3 criteria. The lowest achievable reporting limit for this compound is $0.5~\mu g/L$.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:

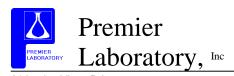
Lynne P. Matteson QA/QC Officer



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

		<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREMEN	TS:	,	
Field measurement calibration records	·	\square	□ _* 11 _*
Groundwater field measurements (if applicable)			NA
Soil sampling field measurements (if applicable)		□ Z	
Sediment sampling field measurements (if applicable)			□ NA
Surface water sampling field measurements (if applicable	:)		אַ(וא □
Low-flow sampling field measurements (if applicable)			□ NA
Documentation of field activities		☑,	
Sample numbering and labeling			
Chain-of-Custody records		I	
Trip blanks		₫,	
Duplicate samples		☑′	مالہ
Equipment blanks			
Split samples (if any)			
2. LABORATORY MEASUREMENTS:	•	,	
Trip blanks		II.	
Instrument blanks			
Laboratory control samples			
Duplicates samples		⊿	
Equipment blanks			□ NA
Matrix spike/matrix spike duplicates		\Box ,	
Analysis type		Ø,	
Chain-of-Custody records		d d	
Surrogate recoveries		I	
Sample Project Narratives		4	· 🗆 👝
Split samples (if any)			□ NA
	1.0	_	_
Т	OTAL: <u>15</u>	<u> </u>	
		1.2	en en
P	ERCENT COM	PLETE:/\	DD %



ANALYTICAL DATA REPORT

Report Number: E711070 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

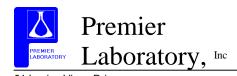
Received Date: 11/1/2007 Report Date: 12/11/2007

> Premier Laboratory, LLC Authorized Signature



Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

Page 1 of 12



		MADEP MCP	Analytical Method I	Report Certificatio	n Form		
Labo	ratory Name: Prei	mier Laboratory, I	nc		Project #: E71	1070	
Proje	ct Location: Frank	klin, MA		l	MADEP RTN ¹ :		
This 1	Form provides cert	ifications for the fo	ollowing data set:[lis	st Laboratory Sam	ple ID Number(s)]	
Samp	ole Matrices: 🛛 🔾	Groundwater 🗵	Soil/Sediment 🗆 🗈	rinking Water	Other		
MCP	SW-846	8260B ⊠	8151A □	8330 🗆	6010B □]	7470A/1A 🗆
Meth	ods Used	8270C □	8081A □	VPH □	6020 🗆		9014M ² □
Compe Analyti (check	ecified in MADEP endium of ical Methods. all that apply)	2 M - SW-846 Metho 3 S - SW-846 Metho	8021B ☐ ing Number (RTN), if kno d 9014 or MADEP Phys ds 7000 Series List indi	iologically Available Cyvidual method and ana	ilyte.	od	7196A □
			ns A, B, C, and D i	-			
Α			aboratory in a condit		า 🛚 🗎	Yes	□ No ¹
_			tody documentation		(-) N		1
В	•	•	ed for the specified a uding the requirement	•	(S) X	Yes	□ No ¹
		•	did not meet appro				
	standards or guid		did flot ffleet approp	onate periornano	=		
С			this report meet all	the requirements	V	Yes	□ No ¹
	•		cribed in Section 2.0	•	ł	100	
		-	ality Assurance and	. , . , . ,			
	for the Acquisition		•				
D			s the VPH or EPH r	nethod run withou	t 🗆	Yes	□ No ¹
	significant modific	cations, as specifie	ed in Section 11.3?				
	A response	to questions E a	nd F below is requ	ired for "Presum	ptive Certainty	r" stat	us
E	Were all QC perfo	ormance standard	s and recommenda	tions for the	X	Yes	□ No ¹
	specified methods	s achieved?					
F	1	•	pounds/elements for	or the specified	X	Yes	□ No ¹
method(s) reported?							
¹ All NO answers must be addressed in an attached Environmental Laboratory case narrative.							
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.							
Sign	Signature: Position: Laboratory Director						
Print	Printed Name: Robert Stevenson Date: 12/11/2007						

Page 2 of 12

> Report No: E711070 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory, Inc received two samples from Fuss & O'Neill on 11/01/2007. The samples were analyzed from the following list of analyses:

Moisture, Percent

Volatiles by 8260B (MCP) in GW/SW 8260B

Variances:

SDG:

None reported.

Method:

None reported.

QA/QC:

None reported.

Page 3 of 12

Laboratory: Premier Laboratory, Inc

PL Report No: E711070 PL Sample No: 1

Date Collected: 10/31/2007
Date Received: 11/1/2007
Date Extracted: By:
Date Analyzed: 11/05/07 By: DDD

Method: 8260B QC Batch#: 57581 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071031-01

Matrix: Solid

Percent Moisture: 9.3 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32638.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	3.2
71-43-2	Benzene	ND	3.2
108-86-1	Bromobenzene	ND	3.2
74-97-5	Bromochloromethane	ND	3.2
75-27-4	Bromodichloromethane	ND	3.2
75-25-2	Bromoform	ND	3.2
74-83-9	Bromomethane	ND	3.2
78-93-3	2-Butanone (MEK)	ND	3.2
104-51-8	n-Butylbenzene	ND	3.2
135-98-8	sec-Butylbenzene	ND	3.2
98-06-6	tert-Butylbenzene	ND	3.2
75-15-0	Carbon disulfide	ND	3.2
56-23-5	Carbon tetrachloride	ND	3.2
108-90-7	Chlorobenzene	ND	3.2
75-00-3	Chloroethane	ND	3.2
67-66-3	Chloroform	ND	3.2
74-87-3	Chloromethane	ND	3.2
95-49-8	2-Chlorotoluene	ND	3.2
106-43-4	4-Chlorotoluene	ND	3.2
108-20-3	Di-isopropyl ether (DIPE)	ND	3.2
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	3.2
124-48-1	Dibromochloromethane	ND	3.2
106-93-4	1,2-Dibromoethane (EDB)	ND	3.2
74-95-3	Dibromomethane	ND	3.2
95-50-1	1,2-Dichlorobenzene	ND	3.2
541-73-1	1,3-Dichlorobenzene	ND	3.2
106-46-7	1,4-Dichlorobenzene	ND	3.2
75-71-8	Dichlorodifluoromethane	ND	3.2
75-34-3	1,1-Dichloroethane	ND	3.2
107-06-2	1,2-Dichloroethane	ND	3.2
75-35-4	1,1-Dichloroethene	ND	3.2
156-59-2	cis-1,2-Dichloroethene	ND	3.2
156-60-5	trans-1,2-Dichloroethene	ND	3.2
78-87-5	1,2-Dichloropropane	ND	3.2
142-28-9	1,3-Dichloropropane	ND	3.2
590-20-7	2,2-Dichloropropane	ND	3.2
563-58-6	1,1-Dichloropropene	ND	3.2
10061-01-5	cis-1,3-Dichloropropene	ND	3.2
10061-02-6	trans-1,3-Dichloropropene	ND	3.2
60-29-7	Diethyl ether	ND	3.2
123-91-1	1,4-Dioxane	ND ND	13
120 /1 1	1,1 DIONUIC	1112	1.5

Page 4 of 12

Laboratory: Premier Laboratory, Inc

Custor

PL Report No: E711070 PL Sample No: 1 (continued)

Date Collected: 10/31/2007
Date Received: 11/1/2007
Date Extracted: By:
Date Analyzed: 11/05/07 By: DDD

Method: 8260B QC Batch#: 57581 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071031-01

Matrix: Solid

Percent Moisture: 9.3 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32638.D

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	3.2
100-41-4	Ethylbenzene	ND	3.2
87-68-3	Hexachlorobutadiene	ND	3.2
591-78-6	2-Hexanone	ND	3.2
98-82-8	Isopropylbenzene	ND	3.2
99-87-6	4-Isopropyltoluene	ND	3.2
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	3.2
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	3.2
75-09-2	Methylene chloride	ND	3.2
91-20-3	Naphthalene	ND	3.2
103-65-1	n-Propylbenzene	ND	3.2
100-42-5	Styrene	ND	3.2
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	3.2
109-99-9	Tetrahydrofuran	ND	3.2
96-18-4	1,2,3-Trichloropropane	ND	3.2
630-20-6	1,1,1,2-Tetrachloroethane	ND	3.2
79-34-5	1,1,2,2-Tetrachloroethane	ND	3.2
127-18-4	Tetrachloroethene (PCE)	6.4	3.2
108-88-3	Toluene	ND	3.2
87-61-6	1,2,3-Trichlorobenzene	ND	3.2
120-82-1	1,2,4-Trichlorobenzene	ND	3.2
71-55-6	1,1,1-Trichloroethane	ND	3.2
79-00-5	1,1,2-Trichloroethane	ND	3.2
79-01-6	Trichloroethene (TCE)	3.5	3.2
75-69-4	Trichlorofluoromethane	ND	3.2
95-63-6	1,2,4-Trimethylbenzene	ND	3.2
108-67-8	1,3,5-Trimethylbenzene	ND	3.2
75-01-4	Vinyl chloride	ND	3.2
95-47-6	o-Xylene	ND	3.2
108-38-3	m,p-Xylenes	ND	3.2
Surrogate	Recovery Lin	nits	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	106%	85%-116%
Bromofluorobenzene	91%	63%-113%
Toluene-d8	90%	78%-128%

Page 5 of 12

Laboratory: Premier Laboratory, Inc

PL Report No: E711070 PL Sample No: 2

Date Collected: 10/31/2007
Date Received: 11/1/2007
Date Extracted: By:
Date Analyzed: 11/05/07 By: DDD

Method: 8260B QC Batch#: 57581 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071031-02

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 50 Soil Extract Volume: Lab Data File: J32637.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	250
71-43-2	Benzene	ND	250
108-86-1	Bromobenzene	ND	250
74-97-5	Bromochloromethane	ND	250
75-27-4	Bromodichloromethane	ND	250
75-25-2	Bromoform	ND	250
74-83-9	Bromomethane	ND	100
78-93-3	2-Butanone (MEK)	ND	250
104-51-8	n-Butylbenzene	ND	250
135-98-8	sec-Butylbenzene	ND	250
98-06-6	tert-Butylbenzene	ND	250
75-15-0	Carbon disulfide	ND	250
56-23-5	Carbon tetrachloride	ND	250
108-90-7	Chlorobenzene	ND	250
75-00-3	Chloroethane	ND	250
67-66-3	Chloroform	ND	250
74-87-3	Chloromethane	ND	250
95-49-8	2-Chlorotoluene	ND	250
106-43-4	4-Chlorotoluene	ND	250
108-20-3	Di-isopropyl ether (DIPE)	ND	250
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	250
124-48-1	Dibromochloromethane	ND	250
106-93-4	1,2-Dibromoethane (EDB)	ND	25
74-95-3	Dibromomethane	ND	250
95-50-1	1,2-Dichlorobenzene	ND	250
541-73-1	1,3-Dichlorobenzene	ND	250
106-46-7	1,4-Dichlorobenzene	ND	250
75-71-8	Dichlorodifluoromethane	ND	250
75-34-3	1,1-Dichloroethane	ND	250
107-06-2	1,2-Dichloroethane	ND	250
75-35-4	1,1-Dichloroethene	ND	50
156-59-2	cis-1,2-Dichloroethene	ND	250
156-60-5	trans-1,2-Dichloroethene	ND	250
78-87-5	1,2-Dichloropropane	ND	250
142-28-9	1,3-Dichloropropane	ND	250
590-20-7	2,2-Dichloropropane	ND	250
563-58-6	1,1-Dichloropropene	ND	250
10061-01-5	cis-1,3-Dichloropropene	ND	25
10061-02-6	trans-1,3-Dichloropropene	ND	25
60-29-7	Diethyl ether	ND	250
123-91-1	1,4-Dioxane	ND	1000

Page 6 of 12

Laboratory: Premier Laboratory, Inc

PL Report No: E711070 PL Sample No: 2 (continued)

Date Collected: 10/31/2007
Date Received: 11/1/2007
Date Extracted: By:
Date Analyzed: 11/05/07 By: DDD

Method: 8260B QC Batch#: 57581 Units: ug/L Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071031-02

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 50 Soil Extract Volume: Lab Data File: J32637.D

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	250
100-41-4	Ethylbenzene	ND	250
87-68-3	Hexachlorobutadiene	ND	30
591-78-6	2-Hexanone	ND	250
98-82-8	Isopropylbenzene	ND	250
99-87-6	4-Isopropyltoluene	ND	250
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	250
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	250
75-09-2	Methylene chloride	ND	250
91-20-3	Naphthalene	ND	250
103-65-1	n-Propylbenzene	ND	250
100-42-5	Styrene	ND	250
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	250
109-99-9	Tetrahydrofuran	ND	250
96-18-4	1,2,3-Trichloropropane	ND	250
630-20-6	1,1,1,2-Tetrachloroethane	ND	250
79-34-5	1,1,2,2-Tetrachloroethane	ND	100
127-18-4	Tetrachloroethene (PCE)	ND	250
108-88-3	Toluene	ND	250
87-61-6	1,2,3-Trichlorobenzene	ND	250
120-82-1	1,2,4-Trichlorobenzene	ND	250
71-55-6	1,1,1-Trichloroethane	ND	250
79-00-5	1,1,2-Trichloroethane	ND	250
79-01-6	Trichloroethene (TCE)	ND	250
75-69-4	Trichlorofluoromethane	ND	250
95-63-6	1,2,4-Trimethylbenzene	ND	250
108-67-8	1,3,5-Trimethylbenzene	ND	250
75-01-4	Vinyl chloride	ND	100
95-47-6	o-Xylene	ND	250
108-38-3	m,p-Xylenes	ND	250
Surrogate	Recovery Lin	nits	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	105%	89%-113%
Bromofluorobenzene	90%	83%-107%
Toluene-d8	93%	88%-108%

Page 7 of 12

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711070 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
1,1,1,2-Tetrachloroet	50.00	45.28	90	72-120
1,1,1-Trichloroethane	50.00	42.97	86	77-123
1,1,2,2-Tetrachloroet	50.00	43.72	87	72-120
1,1,2-Trichloroethane	50.00	43.51	87	80-116
1,1-Dichloroethane	50.00	42.24	84	75-115
1,1-Dichloroethene	50.00	44.20	88	73-129
1,1-Dichloropropene	50.00	42.42	85	75-117
1,2,3-Trichlorobenzene	50.00	43.34	87	70-127
1,2,4-Trichlorobenzene	50.00	47.06	94	70-123
1,2,4-Trimethylbenzene	50.00	47.06	94	76-126
1,2-Dibromoethane (EDB)	50.00	44.08	88	80-114
1,2-Dichlorobenzene	50.00	49.32	99	76-121
1,2-Dichloroethane	50.00	43.54	87	75-115
1,2-Dichloropropane	50.00	43.33	87	71-130
1,3,5-Trimethylbenzene	50.00	45.99	92	76-122
1,3-Dichlorobenzene	50.00	48.99	98	79-120
1,3-Dichloropropane	50.00	43.33	87	82-118
1,4-Dichlorobenzene	50.00	49.20	98	77-116
1,4-Dioxane	50.00	55.46	111	70-119
2,2-Dichloropropane	50.00	43.48	87	71-125
2-Butanone (MEK)	50.00	40.25	80	70-128
4-Chlorotoluene	50.00	48.33	97	77-117
4-Isopropyltoluene	50.00	44.97	90	71-124
4-Methyl-2-pentanone	50.00	43.04	86	75-121
Acetone	50.00	40.07	80	70-118
Benzene	50.00	44.34	89	77-118
Bromobenzene	50.00	47.13	94	79-116
Bromochloromethane	50.00	44.69	89	79-122

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711070 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Bromodichloromethane	50.00	44.14	88	82-127
Bromoform	50.00	46.76	94	78-122
Bromomethane	50.00	45.12	90	70-130
Carbon disulfide	50.00	44.35	89	70-130
Carbon tetrachloride	50.00	48.72	97	77-125
Chlorobenzene	50.00	46.20	92	80-118
Chloroform	50.00	42.94	86	80-113
Chloromethane	50.00	46.15	92	70-130
cis-1,2-Dichloroethene	50.00	44.77	90	85-120
cis-1,3-Dichloropropene	50.00	42.17	84	79-116
Di-isopropyl ether (D	50.00	40.96	82	78-121
Dibromochloromethane	50.00	41.85	84	79-122
Dibromomethane	50.00	45.34	91	78-120
Ethyl tertiary-butyl	50.00	41.23	82	81-122
Ethylbenzene	50.00	49.99	100	84-123
Hexachlorobutadiene	50.00	50.08	100	70-121
Isopropylbenzene	50.00	45.45	91	78-120
m,p-Xylenes	100.0	105.2	105	75-129
Methyl tert-butyl eth	50.00	39.68	79	70-127
Methylene chloride	50.00	40.39	81	72-128
n-Butylbenzene	50.00	44.16	88	70-124
n-Propylbenzene	50.00	47.87	96	80-127
Naphthalene	50.00	43.09	86	70-126
o-Xylene	50.00	48.68	97	78-118
sec-Butylbenzene	50.00	45.08	90	78-118
Styrene	50.00	50.12	100	83-124
tert-Butylbenzene	50.00	46.10	92	76-118
Tertiary-amyl methyl	50.00	42.94	86	81-123

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711070 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Tetrachloroethene (PCE)	50.00	43.97	88	77-122
Toluene	50.00	44.11	88	78-120
trans-1,2-Dichloroethene	50.00	43.65	87	80-120
trans-1,3-Dichloropro	50.00	43.41	87	71-111
Trichloroethene (TCE)	50.00	45.14	90	74-119
Vinyl chloride	50.00	40.70	81	70-116

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711070 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS624 Location: Franklin, MA

Lab File ID: J32630.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
1,1,1-Trichloroethane	20.00	20.70	103	75-125
1,1,2,2-Tetrachloroet	20.00	20.57	103	60-140
1,1,2-Trichloroethane	20.00	19.81	99	71-129
1,1-Dichloroethane	20.00	19.58	98	72-128
1,1-Dichloroethene	20.00	22.89	114	50-150
1,2-Dichlorobenzene	20.00	22.91	114	63-137
1,2-Dichloroethane	20.00	19.68	98	68-132
1,2-Dichloropropane	20.00	20.51	102	34-166
1,3-Dichlorobenzene	20.00	23.68	118	73-127
1,4-Dichlorobenzene	20.00	23.36	117	63-137
2-Chloroethyl vinyl e	20.00	23.24	116	1-224
Benzene	20.00	21.43	107	64-136
Bromodichloromethane	20.00	19.60	98	66-135
Bromoform	20.00	19.22	96	71-129
Bromomethane	20.00	22.70	114	14-186
Carbon tetrachloride	20.00	23.40	117	73-127
Chlorobenzene	20.00	22.36	112	66-134
Chloroethane	20.00	16.71	84	38-162
Chloroform	20.00	19.74	99	68-133
Chloromethane	20.00	17.64	88	1-204
cis-1,3-Dichloropropene	20.00	19.17	96	24-176
Dibromochloromethane	20.00	17.37	87	68-133
Ethylbenzene	20.00	23.80	119	59-141
Methylene chloride	20.00	14.34	72	60-140
Tetrachloroethene (PCE)	20.00	22.36	112	74-127
Toluene	20.00	21.72	108	74-126
trans-1,2-Dichloroethene	20.00	21.40	107	70-131
trans-1,3-Dichloropro	20.00	19.22	96	50-150

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711070 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS624 Location: Franklin, MA

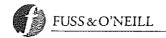
Lab File ID: J32630.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Trichloroethene (TCE)	20.00	21.66	108	66-134
Trichlorofluoromethane	20.00	21.00	105	48-152
Vinyl chloride	20.00	21.41	107	4-196

Column to be used to flag recovery values with an asterisk

* Values outside of QC limits

Page 12 of 12



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS

				<u>YES</u>	NO C	OMMENTS
1.	S	DG Project Narratives		١		
2.		raffic Report			NIT	}
3.	V	olatiles Data				
	a.	<u>t</u>				
		Target Compound List (TCL) Results	M			
		Reconstructed total ion chromatograms (RIC) for each				
		Sample	(2)			
		For each sample:				
		Raw spectra and background-subtracted mass spectra of				
		target compounds identified	A			
		Mass spectra of all reported TICs with three best library matches				
		Percent solids calculations			MIA	
		recent solids calculations	3			
	b.	Standards Data (all instruments)				
		Initial Calibration Data	₽			
		RICs and Quan Reports for all Standards	₩~			
		Continuing Calibration	E			
		RICs and Quan Reports for all Standards				
		Internal Standard Area Summary				
	c.	Raw QC Data				
		Blank Data	\square			
		Matrix Spike Data			MIA	
		Matrix Spike Duplicate Data			NIA	
4.	Sem	nivolatiles Data				
	a. (QC Summary				
		Surrogate Percent Recovery Summary			NIA	
	:	MS/MSD Summary			WIA	
	,	Method Blank Summary			MA	
		Funing and Mass Calibration			MIA	



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

		YES NO	COMMENT
b.	Sample Data		
	TCL Results		NIA
	Tentatively Identified Compounds		NIA
	Reconstructed total ion chromatograms (RIC) for each Sample		NIA
	For each sample:	_	
	Raw spectra and background-subtracted mass		NIA
	spectra of TCL compounds		
	Mass spectra of TICs with 3 best library matches		NIA
	GPC chromatograms (if GPC performed)		WIA
c.	Standards Data (all instruments)		
	Initial Calibration Data		<u> </u>
	RICs and Quan Reports for all Standards		MIA
	Continuing Calibration		W/A
	RICs and Quan Reports for all Standards		N/A
	Internal Standard Areas Summary		NA
	Internal Standard Areas Summary		NIA
d.	Raw QC Data		1 4
	Decafluorotripbenylphosphine (DFTPP)		MIA
	Blank Data		MIA
	Matrix Spike Data		WIA
	Matrix Spike Duplicate Data		WIA
Misc	cellaneous Data		
(Original preparation and analysis forms or copies of preparation		
	and analysis log book pages	Q	
I	Internal sample & sample extract transfer chain-of custody		
	records	№ □	
	Screening Records		NIA
	All instrument output, including strip charts from screening civities (describe or list)		



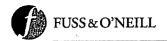
PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

			YES	NO COM	MENTS
6.		ustody Records Sample Log-in Sheet (Lab & DC1) Miscellaneous Shipping/Receiving Records	(describe or list)		4 4
7.	Internal Lab	Sample Transfer Records and Tracking Sheets (describe or list)			
3.	Other Recor	ds (describe or list)			
).	Comments:				
		See laboratory Quality Assurance Plan		rjents Innager	
	Completed (Lab)	by: (Signature) (Prin	regory 1/MK// nted Name/Title)	mager	11/12/07 Date
	I certify that the above as	t the above information is true and accurate nalyses will be made available for review for s	I further certify that all seven (7) years following	laboratory result certification of	s associated with this document.
	Certified by (Lab)		pert Stevenson/Lab X ted Name/Title)	?rector	11-12-67 Date



FUSS & O'NEILL Disciplines to Deliver (860) 646-2469 • www.FandO.com	☐ 146 Hartford Road, Manchester, CT 06040 ☐ 56 Quarry Road, Trumbull, CT 06611 ☐ 1419 Richland Street, Columbia, SC 29201	☐ 610 Lynndale Court, St	est Springfield, MA 01089 nite E, Greenville, NC 27858 ttension, Albany, NY 12203	275 Promenade Street, Suite 35 2 80 Washington Street, Suite 301 2 Other	nade Street, Suite 350, Providence, RI 02908 gton Street, Suite 301, Poughkeepsie, NY 12601				
CHAIN-OF-	CUSTODY RECORD	14718		Turnaroun 1 Day* 3 Days* 2 Days* 4 Standard (days)	d Other (days) *Surcharge Applies				
PROJECT NAME	Project Location		PROJECT NUMBER	Market 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1 1	Laboratory				
Nu-Style Phase II	Franklin, MA		2005045x.Ble	ė .	Premier				
TO: Band Foss		Analysis			Containers				
E To:		Request			/////>				
DE 8410050458 B	010	1	m////////						
's Signature: odes: nitoring Well nice Water PW=Potable Water T=Treatment Facility	Date: 10/31/07 S=Soil W=Waste B=Sediment A=Air	Vox Ob							
Trip Blank		43//							

REPORT	г То:			Sand Foss			7 4		An	alysis		/	///		/	77	/			Co	ontainers	
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Modified Tier II Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E711143
Date Samples Received at Laboratory:	11/2/2007
Date of Review:	12/18/2007

Four soil samples, including a field duplicate, were collected from four soil borings over two days. Samples were submitted to Premier Laboratory in Dayville, Connecticut for analysis of volatile organic compounds (VOCs) by EPA Method 8260B, priority pollutant-13 metals plus barium by EPA Methods 6010B and 7471, and/or petroleum hydrocarbons by Massachusetts Extractable and Volatile Petroleum Hydrocarbon (EPH/VPH) methods of analysis. Analyses for individual samples were determined based upon constituents of concern in the area where the individual monitoring wells are located.

Dedicated sampling equipment was used; therefore, no equipment blank was indicated. An aqueous trip blank was also submitted. No VOCs were reported in the trip blank.

Results of primary and duplicate sample pairs were generally similar. The relative percent difference (RPD) calculated for several metals was greater than the 50% limit established by the QAPP for soil samples. Several RPD values were outside laboratory quality control limits in both initial total and fractionated extractions.

Surrogate recoveries were acceptable for all applicable analyses. With the exception of hexachlorobutadiene, reporting limits were low enough to compare to MADEP GW-2 and GW-3 criteria. The lowest achievable reporting limit for this compound is $0.5~\mu g/L$.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:

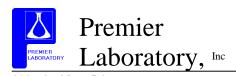
Lynne P. Matteson QA/QC Officer



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

		<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREME	NTS:		
Field measurement calibration records		□ Z	□ , 11,
Groundwater field measurements (if applicable)			
Soil sampling field measurements (if applicable)		√	
Sediment sampling field measurements (if applicable)			□ NIA
Surface water sampling field measurements (if applical	ole)		עומ □
Low-flow sampling field measurements (if applicable)			□ NA
Documentation of field activities		☑,	
Sample numbering and labeling		ব ব ব ব ব	
Chain-of-Custody records		☑	
Trip blanks		\mathbf{Z}_{f}	
Duplicate samples		☑	مالہ 🗆
Equipment blanks			
Split samples (if any)			
2. LABORATORY MEASUREMENTS:		,	
Trip blanks		II,	
Instrument blanks		■ J	
Laboratory control samples		□ ∠	
Duplicates samples		\square	
Equipment blanks			
Matrix spike/matrix spike duplicates			
Analysis type		U	
Chain-of-Custody records			
Surrogate recoveries			
Sample Project Narratives		₫	· □ .0.
Split samples (if any)			□ ~ /A-
	TOTAL:	15	
	PERCENT C	OMPLETE: .	/OD %



ANALYTICAL DATA REPORT

Report Number: E711143 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

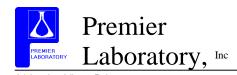
Received Date: 11/2/2007 Report Date: 12/11/2007

> Premier Laboratory, LLC Authorized Signature



Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

Page 1 of 38



MADEP MCP Analytical Method Report Certification Form									
Labor	atory Name: Pre	mier Laboratory, Inc			Project #:	E7111	43		
Proje	ct Location: Fran	klin, MA			MADEP RT	N ¹ :			
This F	orm provides cer	tifications for the follow	ving data set:[list La	boratory Sa	ample ID Num	ber(s)]			
1, 2, 3	3, 4, 5								
Samp	ole Matrices: 🛛 🔾	Groundwater ⊠ Soil/	/Sediment Drinki	ing Water	☐ Other				
MCP	SW-846	8260B ⊠	8151A □	8330	□ 6010	ов ⊠		747	0A/1A ⊠
Meth	ods Used	8270C □	8081A □	VPH [⊠ 602	20 🗆		90)14M ² □
As spe	cified in MADEP	8082 🗆	8021B □	EPH D	⊠ 7000 9	$S^3 \square$		7	7196A 🗆
	endium of	1 List Release Tracking N	umber (RTN), if known		0 :1 (040)				
	cal Methods. all that apply)	2 M - SW-846 Method 901 3 S - SW-846 Methods 70				Method			
-		ponse to questions A				e Certa	aintv		
Α		received by the labora				X Y			No ¹
, ,	•	the Chain-of-Custody	•				00		
В		procedures required for				X Y	'es		No ¹
		port followed, includin		•					
		ative QC data that did	•		nce				
	standards or guid		not moot appropriate	o portorma					
С		al data included in this	report meet all the	requiremen	nts	X Y	'es	$\overline{\Box}$	No ¹
		Certainty", as describe	•	•	1	7 .			
	•	nt CAM VII A, "Quality	, ,	. , . ,	` '				
		n and Reporting of Ana		mry Control					
D	•	ethods only: Was the	•	od run with	out	X Y	'es		No ¹
		cations, as specified in							
		to questions E and F		for "Presi	ımptive Certa	aintv" :	stati		
E		ormance standards an							No ¹
_	specified method			7 101 1110			00	_	
F	•	all analyte-list compou	nds/elements for the	e specified		X Y	'es		No ¹
	method(s) reporte	•		, opeoea					
	` , , ,		d in an attached En	/ironmenta	I I aboratory c	ase na	rrati		
¹ All NO answers must be addressed in an attached Environmental Laboratory case narrative.									
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal									
inquiry of those responsible for obtaining the information, the material contained in this analytical report is, to the best of my knowledge and belief, accurate and complete.									
anaryt	iour report is, to the be	a A	ioi, addarate and demple						
Signa	ature:	Tolor Kurasm	1	Position:	Laboratory Di	rector			
9.1			<u> </u>	30					
Print	ed Name: Robe	rt Stevenson	1	Date: 12/1	1/2007				
	<u> </u>		<u> </u>		00.				

Page 2 of 38

> Report No: E711143 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory, Inc received five samples from Fuss & O'Neill on 11/02/2007. The samples were analyzed from the following list of analyses:

Extractable Petroleum Hydrocarbon (EPH) MADEP EPH[MADEP EPH]

Moisture, Percent

Volatile Petroleum Hydrocarbon (VPH) MADEP VPH

Trace Priority Pollutant (13) Metals in Soil 6010B[3000], 7471[7471] Volatiles by 8260B (MCP) in GW/SW

8260B

Volatiles by 8260B in GW/SW 8260B

Variances:

SDG:

Several RPD values were outside quality control limits for the fractionated LCS. The LCS passed all other quality control limits, in both the initial total and fractionated extractions.

Method:

None reported.

QA/QC:

None reported.

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INORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, Inc PL Report No: E711143

Date Received: 11/2/2007

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(2) 841071101-04					
Date Collected: 11/1/2007 Matrix: Soli	d				
Trace Metals by 6010B	<u>u</u>				
Antimony	1.0	0.19	mg/kg	11/13/07	AMM
Arsenic	3.0	0.32	mg/kg		:06 AMM
Barium	26	0.13	mg/kg		:06 AMM
Beryllium	0.35	0.064	mg/kg		:06 AMM
Cadmium	0.60	0.13	mg/kg		:06 AMM
Chromium	24	0.13	mg/kg		:06 AMM
Copper	110	0.13	mg/kg		:06 AMM
Lead	68	0.13	mg/kg	11/13/07	AMM
Nickel	4.3	0.13	mg/kg		:06 AMM
Selenium	ND	0.32	mg/kg		:06 AMM
Silver	ND	0.13	mg/kg		:06 AMM
Thallium	ND	0.32	mg/kg	11/13/07	AMM
Zinc	73	0.13	mg/kg		:06 AMM
Mercury by SW-846 7471 in SW	0.12	0.026	mg/kg	11/07/07	KAW
(3) 841071101-05					
Date Collected: 11/1/2007 Matrix: Soli	<u>d</u>				
Trace Metals by 6010B					
Antimony	ND	0.18	mg/kg	11/13/07	AMM
Arsenic	1.3	0.30	mg/kg	11/09/07 14	:09 AMM
Barium	15	0.12	mg/kg	11/09/07 14	:09 AMM
Beryllium	0.24	0.061	mg/kg	11/09/07 14	:09 AMM
Cadmium	0.37	0.12	mg/kg	11/09/07 14	:09 AMM
Chromium	15	0.12	mg/kg	11/09/07 14	:09 AMM
Copper	11	0.12	mg/kg	11/09/07 14	:09 AMM
Lead	7.1	0.12	mg/kg	11/13/07	AMM
Nickel	4.7	0.12	mg/kg	11/09/07 14	:09 AMM
Selenium	ND	0.30	mg/kg	11/09/07 14	:09 AMM
Silver	ND	0.12	mg/kg	11/09/07 14	:09 AMM
Thallium	ND	0.30	mg/kg	11/13/07	AMM
Zinc	28	0.12	mg/kg	11/09/07 14	:09 AMM
Mercury by SW-846 7471 in SW	0.028	0.024	mg/kg	11/07/07	KAW

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INORGANIC ANALYSIS DATA SHEET

Laboratory: Premier Laboratory, Inc PL Report No: E711143

Date Received: 11/2/2007

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II

Parameter		Result	DL	Units	Completed		By Dilution
(4) 841071101-06 Date Collected: 11/1/2007	Matrix: Solid						
Trace Metals by 6010B							
Antimony		ND	0.18	mg/kg	11/13/07		AMM
Arsenic		0.84	0.30	mg/kg	11/09/07	14:11	AMM
Barium		15	0.12	mg/kg	11/09/07	14:11	AMM
Beryllium		0.21	0.060	mg/kg	11/09/07	14:11	AMM
Cadmium		0.17	0.12	mg/kg	11/09/07	14:11	AMM
Chromium		3.8	0.12	mg/kg	11/09/07	14:11	AMM
Copper		3.1	0.12	mg/kg	11/09/07	14:11	AMM
Lead		2.5	0.12	mg/kg	11/13/07		AMM
Nickel		2.2	0.12	mg/kg	11/09/07	14:11	AMM
Selenium		ND	0.30	mg/kg	11/09/07	14:11	AMM
Silver		ND	0.12	mg/kg	11/09/07	14:11	AMM
Thallium		ND	0.30	mg/kg	11/13/07		AMM
Zinc		11	0.12	mg/kg	11/09/07	14:11	AMM
Mercury by SW-846 7471 in SW		ND	0.024	mg/kg	11/07/07		KAW

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Laboratory: Premier Laboratory, Inc

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711143 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 Sample Description: 841071101-03

Date Collected: 11/1/2007 Matrix: Solid
Date Received: 11/2/2007 Percent Moiste

Date Received: 11/2/2007 Percent Moisture: 13.7

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 11/06/07 By: DDD Dilution Factor: 200

Method: 8260B Soil Extract Volume:

QC Batch#: 57663 Lab Data File: J32685.D

Units: ug/kg

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	1200
71-43-2	Benzene	ND	1200
108-86-1	Bromobenzene	ND	1200
74-97-5	Bromochloromethane	ND	1200
75-27-4	Bromodichloromethane	ND	1200
75-25-2	Bromoform	ND	1200
74-83-9	Bromomethane	ND	1200
78-93-3	2-Butanone (MEK)	ND	1200
104-51-8	n-Butylbenzene	ND	1200
135-98-8	sec-Butylbenzene	ND	1200
98-06-6	tert-Butylbenzene	ND	1200
75-15-0	Carbon disulfide	ND	1200
56-23-5	Carbon tetrachloride	ND	1200
108-90-7	Chlorobenzene	ND	1200
75-00-3	Chloroethane	ND	1200
67-66-3	Chloroform	ND	1200
74-87-3	Chloromethane	ND	1200
95-49-8	2-Chlorotoluene	ND	1200
106-43-4	4-Chlorotoluene	ND	1200
108-20-3	Di-isopropyl ether (DIPE)	ND	1200
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	1200
124-48-1	Dibromochloromethane	ND	1200
106-93-4	1,2-Dibromoethane (EDB)	ND	1200
74-95-3	Dibromomethane	ND	1200
95-50-1	1,2-Dichlorobenzene	ND	1200
541-73-1	1,3-Dichlorobenzene	ND	1200
106-46-7	1,4-Dichlorobenzene	ND	1200
75-71-8	Dichlorodifluoromethane	ND	1200
75-34-3	1,1-Dichloroethane	ND	1200
107-06-2	1,2-Dichloroethane	ND	1200
75-35-4	1,1-Dichloroethene	ND	1200
156-59-2	cis-1,2-Dichloroethene	ND	1200
156-60-5	trans-1,2-Dichloroethene	ND	1200
78-87-5	1,2-Dichloropropane	ND	1200
142-28-9	1,3-Dichloropropane	ND	1200
590-20-7	2,2-Dichloropropane	ND	1200
563-58-6	1,1-Dichloropropene	ND	1200
10061-01-5	cis-1,3-Dichloropropene	ND	1200
10061-02-6	trans-1,3-Dichloropropene	ND	1200
60-29-7	Diethyl ether	ND	1200
123-91-1	1,4-Dioxane	ND	4600

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Laboratory: Premier Laboratory, Inc

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711143 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 1 (continued) Sample Description: 841071101-03

Date Collected: 11/1/2007 Matrix: Solid

Date Received: 11/2/2007 Percent Moisture: 13.7

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 11/06/07 By: DDD Dilution Factor: 200

Method: 8260B Soil Extract Volume:

QC Batch#: 57663 Lab Data File: J32685.D

Units: ug/kg

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	1200
100-41-4	Ethylbenzene	ND	1200
87-68-3	Hexachlorobutadiene	ND	1200
591-78-6	2-Hexanone	ND	1200
98-82-8	Isopropylbenzene	ND	1200
99-87-6	4-Isopropyltoluene	ND	1200
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1200
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	1200
75-09-2	Methylene chloride	ND	1200
91-20-3	Naphthalene	ND	1200
103-65-1	n-Propylbenzene	ND	1200
100-42-5	Styrene	ND	1200
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	1200
109-99-9	Tetrahydrofuran	ND	1200
96-18-4	1,2,3-Trichloropropane	ND	1200
630-20-6	1,1,1,2-Tetrachloroethane	ND	1200
79-34-5	1,1,2,2-Tetrachloroethane	ND	1200
127-18-4	Tetrachloroethene (PCE)	40000	1200
108-88-3	Toluene	ND	1200
87-61-6	1,2,3-Trichlorobenzene	ND	1200
120-82-1	1,2,4-Trichlorobenzene	ND	1200
71-55-6	1,1,1-Trichloroethane	ND	1200
79-00-5	1,1,2-Trichloroethane	ND	1200
79-01-6	Trichloroethene (TCE)	9200	1200
75-69-4	Trichlorofluoromethane	ND	1200
95-63-6	1,2,4-Trimethylbenzene	ND	1200
108-67-8	1,3,5-Trimethylbenzene	ND	1200
75-01-4	Vinyl chloride	ND	1200
95-47-6	o-Xylene	ND	1200
108-38-3	m,p-Xylenes	ND	1200
Surrogate	Recovery Li	mits	
1.2 Diablamathama 44	020/ 05	20/ 11/0/	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	93%	85%-116%
Bromofluorobenzene	89%	63%-113%
Toluene-d8	93%	78%-128%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, Inc	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711143	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841071101-04
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Analyzed:	11/06/2007	Percent Moisture:	21.9
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6400	ug/kg
C9-C12 Aliphatics**	50	ND	6400	ug/kg
C9-C10 Aromatics***	50	ND	6400	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	80	70%-130%
2,5-dibromotoluene #2	84	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	320	ug/kg
Ethylbenzene	ND	320	ug/kg
Methyl tert-butyl ether (MTBE)	ND	64	ug/kg
Naphthalene	ND	320	ug/kg
Toluene	ND	320	ug/kg
m,p-Xylenes	ND	320	ug/kg
o-Xylene	ND	320	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes *** Excludes Naphthalene

Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 2

Date Collected: 11/1/2007 Date Received: 11/2/2007 Date Extracted: B

Date Analyzed: 11/06/07 By: DDD Method: 8260B QC Batch#: 57663

Units: ug/kg

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-04

Matrix: Solid

Percent Moisture: 21.9 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32680.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	11	6.4
71-43-2	Benzene	ND	6.4
108-86-1	Bromobenzene	ND	6.4
74-97-5	Bromochloromethane	ND	6.4
75-27-4	Bromodichloromethane	ND	6.4
75-25-2	Bromoform	ND	6.4
74-83-9	Bromomethane	ND	6.4
78-93-3	2-Butanone (MEK)	ND	6.4
104-51-8	n-Butylbenzene	ND	6.4
135-98-8	sec-Butylbenzene	ND	6.4
98-06-6	tert-Butylbenzene	ND	6.4
75-15-0	Carbon disulfide	ND	6.4
56-23-5	Carbon tetrachloride	ND	6.4
108-90-7	Chlorobenzene	ND	6.4
75-00-3	Chloroethane	ND	6.4
67-66-3	Chloroform	ND	6.4
74-87-3	Chloromethane	ND	6.4
95-49-8	2-Chlorotoluene	ND	6.4
106-43-4	4-Chlorotoluene	ND	6.4
108-20-3	Di-isopropyl ether (DIPE)	ND	6.4
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	6.4
124-48-1	Dibromochloromethane	ND	6.4
106-93-4	1,2-Dibromoethane (EDB)	ND	6.4
74-95-3	Dibromomethane	ND	6.4
95-50-1	1,2-Dichlorobenzene	ND	6.4
541-73-1	1,3-Dichlorobenzene	ND	6.4
106-46-7	1,4-Dichlorobenzene	ND	6.4
75-71-8	Dichlorodifluoromethane	ND	6.4
75-34-3	1,1-Dichloroethane	ND	6.4
107-06-2	1,2-Dichloroethane	ND	6.4
75-35-4	1,1-Dichloroethene	ND	6.4
156-59-2	cis-1,2-Dichloroethene	ND	6.4
156-60-5	trans-1,2-Dichloroethene	ND	6.4
78-87-5	1,2-Dichloropropane	ND	6.4
142-28-9	1,3-Dichloropropane	ND	6.4
590-20-7	2,2-Dichloropropane	ND	6.4
563-58-6	1,1-Dichloropropene	ND	6.4
10061-01-5	cis-1,3-Dichloropropene	ND	6.4
10061-02-6	trans-1,3-Dichloropropene	ND	6.4
60-29-7	Diethyl ether	ND	6.4
123-91-1	1,4-Dioxane	ND	26

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Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 2 (continued)

Date Collected: 11/1/2007 Date Received: 11/2/2007 Date Extracted: B

Date Analyzed: 11/06/07 By: DDD

Method: 8260B QC Batch#: 57663 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-04

Matrix: Solid

Percent Moisture: 21.9 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32680.D

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	6.4
100-41-4	Ethylbenzene	ND	6.4
87-68-3	Hexachlorobutadiene	ND	6.4
591-78-6	2-Hexanone	ND	6.4
98-82-8	Isopropylbenzene	ND	6.4
99-87-6	4-Isopropyltoluene	ND	6.4
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	6.4
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	6.4
75-09-2	Methylene chloride	ND	6.4
91-20-3	Naphthalene	140	6.4
103-65-1	n-Propylbenzene	ND	6.4
100-42-5	Styrene	ND	6.4
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	6.4
109-99-9	Tetrahydrofuran	ND	6.4
96-18-4	1,2,3-Trichloropropane	ND	6.4
630-20-6	1,1,1,2-Tetrachloroethane	ND	6.4
79-34-5	1,1,2,2-Tetrachloroethane	ND	6.4
127-18-4	Tetrachloroethene (PCE)	ND	6.4
108-88-3	Toluene	ND	6.4
87-61-6	1,2,3-Trichlorobenzene	ND	6.4
120-82-1	1,2,4-Trichlorobenzene	ND	6.4
71-55-6	1,1,1-Trichloroethane	ND	6.4
79-00-5	1,1,2-Trichloroethane	ND	6.4
79-01-6	Trichloroethene (TCE)	ND	6.4
75-69-4	Trichlorofluoromethane	ND	6.4
95-63-6	1,2,4-Trimethylbenzene	ND	6.4
108-67-8	1,3,5-Trimethylbenzene	ND	6.4
75-01-4	Vinyl chloride	ND	6.4
95-47-6	o-Xylene	ND	6.4
108-38-3	m,p-Xylenes	ND	6.4
Surrogate	Recovery Lin	nits	
1.2 Diablamathama 44	1010/ 950	/ 11.00/	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	101%	85%-116%
Bromofluorobenzene	85%	63%-113%
Toluene-d8	100%	78%-128%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, Inc	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711143	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841071101-05
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Analyzed:	11/06/2007	Percent Moisture:	18.0
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6100	ug/kg
C9-C12 Aliphatics**	50	ND	6100	ug/kg
C9-C10 Aromatics***	50	ND	6100	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	80	70%-130%
2,5-dibromotoluene #2	89	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	300	ug/kg
Ethylbenzene	ND	300	ug/kg
Methyl tert-butyl ether (MTBE)	ND	61	ug/kg
Naphthalene	ND	300	ug/kg
Toluene	ND	300	ug/kg
m,p-Xylenes	ND	300	ug/kg
o-Xylene	ND	300	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes *** Excludes Naphthalene

Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 3

Date Collected: 11/1/2007 Date Received: 11/2/2007 Date Extracted: By:

Date Analyzed: 11/05/07 By: DDD Method: 8260B QC Batch#: 57581

Units: ug/kg

CAS No.

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-05

Result

DL

Matrix: Solid

Percent Moisture: 18.0 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32641.D

CAS NO.	Farameter	Kesuit	DL
67-64-1	Acetone	28	4.7
71-43-2	Benzene	ND	4.7
108-86-1	Bromobenzene	ND	4.7
74-97-5	Bromochloromethane	ND	4.7
75-27-4	Bromodichloromethane	ND	4.7
75-25-2	Bromoform	ND	4.7
74-83-9	Bromomethane	ND	4.7
78-93-3	2-Butanone (MEK)	6.6	4.7
104-51-8	n-Butylbenzene	ND	4.7
135-98-8	sec-Butylbenzene	ND	4.7
98-06-6	tert-Butylbenzene	ND	4.7
75-15-0	Carbon disulfide	ND	4.7
56-23-5	Carbon tetrachloride	ND	4.7
108-90-7	Chlorobenzene	ND	4.7
75-00-3	Chloroethane	ND	4.7
67-66-3	Chloroform	ND	4.7
74-87-3	Chloromethane	ND	4.7
95-49-8	2-Chlorotoluene	ND	4.7
106-43-4	4-Chlorotoluene	ND	4.7
108-20-3	Di-isopropyl ether (DIPE)	ND	4.7
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	4.7
124-48-1	Dibromochloromethane	ND	4.7
106-93-4	1,2-Dibromoethane (EDB)	ND	4.7
74-95-3	Dibromomethane	ND	4.7
95-50-1	1,2-Dichlorobenzene	ND	4.7
541-73-1	1,3-Dichlorobenzene	ND	4.7
106-46-7	1,4-Dichlorobenzene	ND	4.7
75-71-8	Dichlorodifluoromethane	ND	4.7
75-34-3	1,1-Dichloroethane	ND	4.7
107-06-2	1,2-Dichloroethane	ND	4.7
75-35-4	1,1-Dichloroethene	ND	4.7
156-59-2	cis-1,2-Dichloroethene	ND	4.7
156-60-5	trans-1,2-Dichloroethene	ND	4.7
78-87-5	1,2-Dichloropropane	ND	4.7
142-28-9	1,3-Dichloropropane	ND	4.7
590-20-7	2,2-Dichloropropane	ND	4.7
563-58-6	1,1-Dichloropropene	ND	4.7
10061-01-5	cis-1,3-Dichloropropene	ND	4.7
10061-02-6	trans-1,3-Dichloropropene	ND	4.7
60-29-7	Diethyl ether	ND	4.7
123-91-1	1,4-Dioxane	ND	19

Parameter

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Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 3 (continued)

Date Collected: 11/1/2007 Date Received: 11/2/2007 Date Extracted: By

Date Analyzed: 11/05/07 By: DDD Method: 8260B

QC Batch#: 57581 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-05

Matrix: Solid

Percent Moisture: 18.0 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32641.D

CAS No.	Parameter	Result	DL
	Ethyl tertiary-butyl ether (EtBE)	ND	4.7
100-41-4	Ethylbenzene	ND	4.7
87-68-3	Hexachlorobutadiene	ND	4.7
591-78-6	2-Hexanone	ND	4.7
98-82-8	Isopropylbenzene	ND	4.7
99-87-6	4-Isopropyltoluene	ND	4.7
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	4.7
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	4.7
75-09-2	Methylene chloride	ND	4.7
91-20-3	Naphthalene	ND	4.7
103-65-1	n-Propylbenzene	ND	4.7
100-42-5	Styrene	ND	4.7
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	4.7
109-99-9	Tetrahydrofuran	ND	4.7
96-18-4	1,2,3-Trichloropropane	ND	4.7
630-20-6	1,1,1,2-Tetrachloroethane	ND	4.7
79-34-5	1,1,2,2-Tetrachloroethane	ND	4.7
127-18-4	Tetrachloroethene (PCE)	ND	4.7
108-88-3	Toluene	ND	4.7
87-61-6	1,2,3-Trichlorobenzene	ND	4.7
120-82-1	1,2,4-Trichlorobenzene	ND	4.7
71-55-6	1,1,1-Trichloroethane	ND	4.7
79-00-5	1,1,2-Trichloroethane	ND	4.7
79-01-6	Trichloroethene (TCE)	ND	4.7
75-69-4	Trichlorofluoromethane	ND	4.7
95-63-6	1,2,4-Trimethylbenzene	ND	4.7
108-67-8	1,3,5-Trimethylbenzene	ND	4.7
75-01-4	Vinyl chloride	ND	4.7
95-47-6	o-Xylene	ND	4.7
108-38-3	m,p-Xylenes	ND	4.7
Surrogate	Recovery Limits		

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	107%	85%-116%
Bromofluorobenzene	83%	63%-113%
Toluene-d8	89%	78%-128%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, Inc	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711143	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	4	Sample Description:	841071101-06
Preservative	METHANOL		
		Dilution (Target):	50
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Analyzed:	11/06/2007	Percent Moisture:	17.2
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	50	ND	6000	ug/kg
C9-C12 Aliphatics**	50	ND	6000	ug/kg
C9-C10 Aromatics***	50	ND	6000	ug/kg

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene	76	70%-130%
2,5-dibromotoluene #2	88	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	QL	Units
Benzene	ND	300	ug/kg
Ethylbenzene	ND	300	ug/kg
Methyl tert-butyl ether (MTBE)	ND	60	ug/kg
Naphthalene	ND	300	ug/kg
Toluene	ND	300	ug/kg
m,p-Xylenes	ND	300	ug/kg
o-Xylene	ND	300	ug/kg

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^{**} Excludes Ethylbenzene, Xylenes *** Excludes Naphthalene

Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 4

Date Collected: 11/1/2007
Date Received: 11/2/2007
Date Extracted: By:

Date Analyzed: 11/05/07 By: DDD Method: 8260B

Units: ug/kg

CAS No.

QC Batch#: 57581

Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-06

Result

DL

Matrix: Solid

Percent Moisture: 17.2 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32642.D

CAS NO.	Farameter	Kesuit	<u>DL</u>
67-64-1	Acetone	35	5.2
71-43-2	Benzene	ND	5.2
108-86-1	Bromobenzene	ND	5.2
74-97-5	Bromochloromethane	ND	5.2
75-27-4	Bromodichloromethane	ND	5.2
75-25-2	Bromoform	ND	5.2
74-83-9	Bromomethane	ND	5.2
78-93-3	2-Butanone (MEK)	7.4	5.2
104-51-8	n-Butylbenzene	ND	5.2
135-98-8	sec-Butylbenzene	ND	5.2
98-06-6	tert-Butylbenzene	ND	5.2
75-15-0	Carbon disulfide	ND	5.2
56-23-5	Carbon tetrachloride	ND	5.2
108-90-7	Chlorobenzene	ND	5.2
75-00-3	Chloroethane	ND	5.2
67-66-3	Chloroform	ND	5.2
74-87-3	Chloromethane	ND	5.2
95-49-8	2-Chlorotoluene	ND	5.2
106-43-4	4-Chlorotoluene	ND	5.2
108-20-3	Di-isopropyl ether (DIPE)	ND	5.2
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	5.2
124-48-1	Dibromochloromethane	ND	5.2
106-93-4	1,2-Dibromoethane (EDB)	ND	5.2
74-95-3	Dibromomethane	ND	5.2
95-50-1	1,2-Dichlorobenzene	ND	5.2
541-73-1	1,3-Dichlorobenzene	ND	5.2
106-46-7	1,4-Dichlorobenzene	ND	5.2
75-71-8	Dichlorodifluoromethane	ND	5.2
75-34-3	1,1-Dichloroethane	ND	5.2
107-06-2	1,2-Dichloroethane	ND	5.2
75-35-4	1,1-Dichloroethene	ND	5.2
156-59-2	cis-1,2-Dichloroethene	ND	5.2
156-60-5	trans-1,2-Dichloroethene	ND	5.2
78-87-5	1,2-Dichloropropane	ND	5.2
142-28-9	1,3-Dichloropropane	ND	5.2
590-20-7	2,2-Dichloropropane	ND	5.2
563-58-6	1,1-Dichloropropene	ND	5.2
10061-01-5	cis-1,3-Dichloropropene	ND	5.2
10061-02-6	trans-1,3-Dichloropropene	ND	5.2
60-29-7	Diethyl ether	ND	5.2
123-91-1	1,4-Dioxane	ND	21

Parameter

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Laboratory: Premier Laboratory, Inc

PL Report No: E711143 PL Sample No: 4 (continued)

Date Collected: 11/1/2007 Date Received: 11/2/2007 Date Extracted: B

Date Analyzed: 11/05/07 By: DDD Method: 8260B

QC Batch#: 57581 Units: ug/kg Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II Sample Description: 841071101-06

Matrix: Solid

Percent Moisture: 17.2 Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32642.D

CAS No.	Parameter		DL
	Ethyl tertiary-butyl ether (EtBE)	ND	5.2
100-41-4	Ethylbenzene	ND	5.2
87-68-3	Hexachlorobutadiene	ND	5.2
591-78-6	2-Hexanone	ND	5.2
98-82-8	Isopropylbenzene	ND	5.2
99-87-6	4-Isopropyltoluene	ND	5.2
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	5.2
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.2
75-09-2	Methylene chloride	ND	5.2
91-20-3	Naphthalene	ND	5.2
103-65-1	n-Propylbenzene	ND	5.2
100-42-5	Styrene	ND	5.2
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	5.2
109-99-9	Tetrahydrofuran	ND	5.2
96-18-4	1,2,3-Trichloropropane	ND	5.2
630-20-6	1,1,1,2-Tetrachloroethane	ND	5.2
79-34-5	1,1,2,2-Tetrachloroethane	ND	5.2
127-18-4	Tetrachloroethene (PCE)	ND	5.2
108-88-3	Toluene	ND	5.2
87-61-6	1,2,3-Trichlorobenzene	ND	5.2
120-82-1	1,2,4-Trichlorobenzene	ND	5.2
71-55-6	1,1,1-Trichloroethane	ND	5.2
79-00-5	1,1,2-Trichloroethane	ND	5.2
79-01-6	Trichloroethene (TCE)	ND	5.2
75-69-4	Trichlorofluoromethane	ND	5.2
95-63-6	1,2,4-Trimethylbenzene	ND	5.2
108-67-8	1,3,5-Trimethylbenzene	ND	5.2
75-01-4	Vinyl chloride	ND	5.2
95-47-6	o-Xylene	ND	5.2
108-38-3	m,p-Xylenes	ND	5.2
Surrogate	Recovery Limits		

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	105%	85%-116%
Bromofluorobenzene	85%	63%-113%
Toluene-d8	90%	78%-128%

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711143 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841071101-07

Date Collected: 11/1/2007 Matrix: Aqueous
Date Received: 11/2/2007 Percent Moisture: N/A
Date Extracted: By: Sample Weight/Volume:
Date Analyzed: 11/05/07 By: DDD Dilution Factor: 50
Mathed: 8260B

Method: 8260B Soil Extract Volume:
QC Batch#: 57581 Lab Data File: J32636.D
Units: ug/L

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	250
71-43-2	Benzene	ND	250
108-86-1	Bromobenzene	ND	250
74-97-5	Bromochloromethane	ND	250
75-27-4	Bromodichloromethane	ND	250
75-25-2	Bromoform	ND	250
74-83-9	Bromomethane	ND	100
78-93-3	2-Butanone (MEK)	ND	250
104-51-8	n-Butylbenzene	ND	250
135-98-8	sec-Butylbenzene	ND	250
98-06-6	tert-Butylbenzene	ND	250
75-15-0	Carbon disulfide	ND	250
56-23-5	Carbon tetrachloride	ND	250
108-90-7	Chlorobenzene	ND	250
75-00-3	Chloroethane	ND	250
67-66-3	Chloroform	ND	250
74-87-3	Chloromethane	ND	250
95-49-8	2-Chlorotoluene	ND	250
106-43-4	4-Chlorotoluene	ND	250
108-20-3	Di-isopropyl ether (DIPE)	ND	250
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	250
124-48-1	Dibromochloromethane	ND	250
106-93-4	1,2-Dibromoethane (EDB)	ND	25
74-95-3	Dibromomethane	ND	250
95-50-1	1,2-Dichlorobenzene	ND	250
541-73-1	1,3-Dichlorobenzene	ND	250
106-46-7	1,4-Dichlorobenzene	ND	250
75-71-8	Dichlorodifluoromethane	ND	250
75-34-3	1,1-Dichloroethane	ND	250
107-06-2	1,2-Dichloroethane	ND	250
75-35-4	1,1-Dichloroethene	ND	50
156-59-2	cis-1,2-Dichloroethene	ND	250
156-60-5	trans-1,2-Dichloroethene	ND	250
78-87-5	1,2-Dichloropropane	ND	250
142-28-9	1,3-Dichloropropane	ND	250
590-20-7	2,2-Dichloropropane	ND	250
563-58-6	1,1-Dichloropropene	ND	250
10061-01-5	cis-1,3-Dichloropropene	ND	25
10061-02-6	trans-1,3-Dichloropropene	ND	25
60-29-7	Diethyl ether	ND	250
123-91-1	1,4-Dioxane	ND	1000

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Lab Data File: J32636.D

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711143 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 (continued) Sample Description: 841071101-07

Date Collected: 11/1/2007 Matrix: Aqueous
Date Received: 11/2/2007 Percent Moisture: N/A
Date Extracted: By: Sample Weight/Volume:
Date Analyzed: 11/05/07 By: DDD Dilution Factor: 50
Method: 8260B Soil Extract Volume:

QC Batch#: 57581 Units: ug/L

Toluene-d8

CAS No.	No. Parameter		DL
	Ethyl tertiary-butyl ether (EtBE)	ND	250
100-41-4	Ethylbenzene	ND	250
87-68-3	Hexachlorobutadiene	ND	30
591-78-6	2-Hexanone	ND	250
98-82-8	Isopropylbenzene	ND	250
99-87-6	4-Isopropyltoluene	ND	250
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	250
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	250
75-09-2	Methylene chloride	ND	250
91-20-3	Naphthalene	ND	250
103-65-1	n-Propylbenzene	ND	250
100-42-5	Styrene	ND	250
994-05-8	Tertiary-amyl methyl ether (TAME)	ND	250
109-99-9	Tetrahydrofuran	ND	250
96-18-4	1,2,3-Trichloropropane	ND	250
630-20-6	1,1,1,2-Tetrachloroethane	ND	250
79-34-5	1,1,2,2-Tetrachloroethane	ND	100
127-18-4	Tetrachloroethene (PCE)	ND	250
108-88-3	Toluene		
87-61-6	1,2,3-Trichlorobenzene		
120-82-1	1,2,4-Trichlorobenzene		
71-55-6	1,1,1-Trichloroethane	ND	250
79-00-5	1,1,2-Trichloroethane	ND	250
79-01-6	Trichloroethene (TCE)	ND	250
75-69-4	Trichlorofluoromethane	ND	250
95-63-6	1,2,4-Trimethylbenzene	ND	250
108-67-8	1,3,5-Trimethylbenzene	ND	250
75-01-4	Vinyl chloride	ND	100
95-47-6	o-Xylene	ND	250
108-38-3	m,p-Xylenes	ND	250
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	104%	81%-115%	
Bromofluorobenzene	90%	85%-106%	

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88%

83%-114%

EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, Inc	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711143	Project:	20050458.B10/Nu-Style Phase II
	_, _, _,	J	841071101-04
PL Sample No:	2	Sample Description:	841071101-04
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Extracted:	11/06/2007	Percent Moisture:	21.9
Date Analyzed:	11/07/2007	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	13000	ug/kg
C19-C36 Aliphatics	1	18000	13000	ug/kg
C11-C22 Aromatics*	1	60000	13000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	62	40%-140%
2-Bromonaphthalene	67	40%-140%
2-Fluorobiphenyl	68	40%-140%
o-Terphenyl	48	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	130	ug/kg
Acenaphthene	150	130	ug/kg
Acenaphthylene	260	130	ug/kg
Anthracene	640	130	ug/kg
Benzo[a]anthracene	1800	130	ug/kg
Benzo[a]pyrene	1600	130	ug/kg
Benzo[b]fluoranthene	1800	130	ug/kg
Benzo[g,h,i]perylene	ND	130	ug/kg
Benzo[k]fluoranthene	1300	130	ug/kg
Chrysene	1700	130	ug/kg
Dibenz[a,h]anthracene	ND	130	ug/kg
Fluoranthene	3900	130	ug/kg
Fluorene	280	130	ug/kg
Indeno[1,2,3-cd]pyrene	1200	130	ug/kg
Naphthalene	130	130	ug/kg
Phenanthrene	2800	130	ug/kg
Pyrene	3600	130	ug/kg

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, Inc	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711143	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841071101-05
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Extracted:	11/06/2007	Percent Moisture:	18.0
Date Analyzed:	11/07/2007	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	12000	ug/kg
C19-C36 Aliphatics	1	ND	12000	ug/kg
C11-C22 Aromatics*	1	17000	12000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	49	40%-140%
2-Bromonaphthalene	83	40%-140%
2-Fluorobiphenyl	80	40%-140%
o-Terphenyl	61	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	120	ug/kg
Acenaphthene	ND	120	ug/kg
Acenaphthylene	ND	120	ug/kg
Anthracene	ND	120	ug/kg
Benzo[a]anthracene	ND	120	ug/kg
Benzo[a]pyrene	ND	120	ug/kg
Benzo[b]fluoranthene	ND	120	ug/kg
Benzo[g,h,i]perylene	ND	120	ug/kg
Benzo[k]fluoranthene	ND	120	ug/kg
Chrysene	ND	120	ug/kg
Dibenz[a,h]anthracene	ND	120	ug/kg
Fluoranthene	180	120	ug/kg
Fluorene	ND	120	ug/kg
Indeno[1,2,3-cd]pyrene	ND	120	ug/kg
Naphthalene	ND	120	ug/kg
Phenanthrene	290	120	ug/kg
Pyrene	180	120	ug/kg

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, Inc	Client:	Fuss & O'Neill
DI Damout No.	E711143	Location:	Franklin, MA 20050458.B10/Nu-Style Phase II
PL Report No:	E/11143	Project:	•
PL Sample No:	4	Sample Description:	841071101-06
Preservative	None		
		Dilution (Target):	1
Date Collected:	11/1/2007		
Date Received:	11/2/2007	Matrix:	Solid
Date Extracted:	11/06/2007	Percent Moisture:	17.2
Date Analyzed:	11/07/2007	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	12000	ug/kg
C19-C36 Aliphatics	1	ND	12000	ug/kg
C11-C22 Aromatics*	1	ND	12000	ug/kg

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	69	40%-140%
2-Bromonaphthalene	59	40%-140%
2-Fluorobiphenyl	59	40%-140%
o-Terphenyl	96	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	120	ug/kg
Acenaphthene	ND	120	ug/kg
Acenaphthylene	ND	120	ug/kg
Anthracene	ND	120	ug/kg
Benzo[a]anthracene	ND	120	ug/kg
Benzo[a]pyrene	ND	120	ug/kg
Benzo[b]fluoranthene	ND	120	ug/kg
Benzo[g,h,i]perylene	ND	120	ug/kg
Benzo[k]fluoranthene	ND	120	ug/kg
Chrysene	ND	120	ug/kg
Dibenz[a,h]anthracene	ND	120	ug/kg
Fluoranthene	ND	120	ug/kg
Fluorene	ND	120	ug/kg
Indeno[1,2,3-cd]pyrene	ND	120	ug/kg
Naphthalene	ND	120	ug/kg
Phenanthrene	ND	120	ug/kg
Pyrene	ND	120	ug/kg

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FORM 3 Soil 6010B Lab Control Sample

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/09/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS Location: Franklin, MA

Lab File ID: 110907A.lim

	Spike Added	Sample Concentration	%	QC Limits
Compound	(ppb)	(ppb)	Rec#	Rec
Aluminum	(PP-07	(PP-0)		-
Antimony				_
Arsenic				_
Barium				-
Beryllium				-
Boron				-
Cadmium				-
Calcium				-
Chromium				-
Cobalt				-
Copper				-
Iron				-
Lead				-
Magnesium				-
Manganese				-
Molybdenum				-
Nickel				-
Potassium				-
Selenium				-
Silver				-
Sodium				-
Thallium				-
Tin				-
Titanium				-
Vanadium				-
Zinc				_

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 4 6010B Method Blank Summary

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Lab File ID: 110907A.lim Lab Sample ID: Blank

Matrix: Soil Date Analyzed: 11/09/07

Instrument ID: VarICP Date Extracted:

Time Analyzed: 1354

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	LCS	LCS	110907A.lim	Analyzed 11/09/2007
1 2 3				
3				
4				
5				
4 5 6				
7				
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8 9				
10				
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Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
1,1,1,2-Tetrachloroet	50.00	45.28	90	72-120
1,1,1-Trichloroethane	50.00	42.97	86	77-123
1,1,2,2-Tetrachloroet	50.00	43.72	87	72-120
1,1,2-Trichloroethane	50.00	43.51	87	80-116
1,1-Dichloroethane	50.00	42.24	84	75-115
1,1-Dichloroethene	50.00	44.20	88	73-129
1,1-Dichloropropene	50.00	42.42	85	75-117
1,2,3-Trichlorobenzene	50.00	43.34	87	70-127
1,2,4-Trichlorobenzene	50.00	47.06	94	70-123
1,2,4-Trimethylbenzene	50.00	47.06	94	76-126
1,2-Dibromoethane (EDB)	50.00	44.08	88	80-114
1,2-Dichlorobenzene	50.00	49.32	99	76-121
1,2-Dichloroethane	50.00	43.54	87	75-115
1,2-Dichloropropane	50.00	43.33	87	71-130
1,3,5-Trimethylbenzene	50.00	45.99	92	76-122
1,3-Dichlorobenzene	50.00	48.99	98	79-120
1,3-Dichloropropane	50.00	43.33	87	82-118
1,4-Dichlorobenzene	50.00	49.20	98	77-116
1,4-Dioxane	50.00	55.46	111	70-119
2,2-Dichloropropane	50.00	43.48	87	71-125
2-Butanone (MEK)	50.00	40.25	80	70-128
4-Chlorotoluene	50.00	48.33	97	77-117
4-Isopropyltoluene	50.00	44.97	90	71-124
4-Methyl-2-pentanone	50.00	43.04	86	75-121
Acetone	50.00	40.07	80	70-118
Benzene	50.00	44.34	89	77-118
Bromobenzene	50.00	47.13	94	79-116
Bromochloromethane	50.00	44.69	89	79-122

[#] Column to be used to flag recovery values with an asterisk

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^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Bromodichloromethane	50.00	44.14	88	82-127
Bromoform	50.00	46.76	94	78-122
Bromomethane	50.00	45.12	90	70-130
Carbon disulfide	50.00	44.35		70-130
Carbon tetrachloride	50.00	48.72	97	77-125
Chlorobenzene	50.00	46.20	92	80-118
Chloroform	50.00	42.94	86	80-113
Chloromethane	50.00	46.15	92	70-130
cis-1,2-Dichloroethene	50.00	44.77	90	85-120
cis-1,3-Dichloropropene	50.00	42.17	84	79-116
Di-isopropyl ether (D	50.00	40.96	82	78-121
Dibromochloromethane	50.00	41.85	84	79-122
Dibromomethane	50.00	45.34	91	78-120
Ethyl tertiary-butyl	50.00	41.23	82	81-122
Ethylbenzene	50.00	49.99	100	84-123
Hexachlorobutadiene	50.00	50.08	100	70-121
Isopropylbenzene	50.00	45.45	91	78-120
m,p-Xylenes	100.0	105.2		75-129
Methyl tert-butyl eth	50.00	39.68	79	70-127
Methylene chloride	50.00	40.39	81	72-128
n-Butylbenzene	50.00	44.16	88	70-124
n-Propylbenzene	50.00	47.87	96	80-127
Naphthalene	50.00	43.09	86	70-126
o-Xylene	50.00	48.68	97	78-118
sec-Butylbenzene	50.00	45.08	90	78-118
Styrene	50.00	50.12	100	83-124
tert-Butylbenzene	50.00	46.10	92	76-118
Tertiary-amyl methyl	50.00	42.94	86	81-123

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1105 Location: Franklin, MA

Lab File ID: J32628.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Tetrachloroethene (PCE)	50.00	43.97	88	77-122
Toluene	50.00	44.11	88	78-120
trans-1,2-Dichloroethene	50.00	43.65	87	80-120
trans-1,3-Dichloropro	50.00	43.41	87	71-111
Trichloroethene (TCE)	50.00	45.14	90	74-119
Vinyl chloride	50.00	40.70	81	70-116

[#] Column to be used to flag recovery values with an asterisk

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^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS624 Location: Franklin, MA

Lab File ID: J32630.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
1,1,1-Trichloroethane	20.00	20.70	103	75-125
1,1,2,2-Tetrachloroet	20.00	20.57	103	60-140
1,1,2-Trichloroethane	20.00	19.81	99	71-129
1,1-Dichloroethane	20.00	19.58	98	72-128
1,1-Dichloroethene	20.00	22.89	114	50-150
1,2-Dichlorobenzene	20.00	22.91	114	63-137
1,2-Dichloroethane	20.00	19.68	98	68-132
1,2-Dichloropropane	20.00	20.51	102	34-166
1,3-Dichlorobenzene	20.00	23.68	118	73-127
1,4-Dichlorobenzene	20.00	23.36	117	63-137
2-Chloroethyl vinyl e	20.00	23.24	116	1-224
Benzene	20.00	21.43	107	64-136
Bromodichloromethane	20.00	19.60	98	66-135
Bromoform	20.00	19.22	96	71-129
Bromomethane	20.00	22.70	114	14-186
Carbon tetrachloride	20.00	23.40	117	73-127
Chlorobenzene	20.00	22.36	112	66-134
Chloroethane	20.00	16.71	84	38-162
Chloroform	20.00	19.74	99	68-133
Chloromethane	20.00	17.64	88	1-204
cis-1,3-Dichloropropene	20.00	19.17	96	24-176
Dibromochloromethane	20.00	17.37	87	68-133
Ethylbenzene	20.00	23.80	119	59-141
Methylene chloride	20.00	14.34	72	60-140
Tetrachloroethene (PCE)	20.00	22.36	112	74-127
Toluene	20.00	21.72	108	74-126
trans-1,2-Dichloroethene	20.00	21.40	107	70-131
trans-1,3-Dichloropro	20.00	19.22	96	50-150

[#] Column to be used to flag recovery values with an asterisk

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^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/05/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS624 Location: Franklin, MA

Lab File ID: J32630.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Trichloroethene (TCE)	20.00	21.66	108	66-134
Trichlorofluoromethane	20.00	21.00	105	48-152
Vinyl chloride	20.00	21.41	107	4-196

Column to be used to flag recovery values with an asterisk

* Values outside of QC limits

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Lab Name: Premier Laboratory, Inc Date Analyzed: 11/06/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1106 Location: Franklin, MA

Lab File ID: J32674B.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
1,1,1,2-Tetrachloroet	50.00	45.93	92	72-120
1,1,1-Trichloroethane	50.00	45.83	92	77-123
1,1,2,2-Tetrachloroet	50.00	50.44	101	72-120
1,1,2-Trichloroethane	50.00	45.97	92	80-116
1,1-Dichloroethane	50.00	49.35	99	75-115
1,1-Dichloroethene	50.00	54.20	108	73-129
1,1-Dichloropropene	50.00	47.98	96	75-117
1,2,3-Trichlorobenzene	50.00	45.95	92	70-127
1,2,4-Trichlorobenzene	50.00	45.37	91	70-123
1,2,4-Trimethylbenzene	50.00	42.74	85	76-126
1,2-Dibromoethane (EDB)	50.00	45.70	91	80-114
1,2-Dichlorobenzene	50.00	52.61	105	76-121
1,2-Dichloroethane	50.00	47.82	96	75-115
1,2-Dichloropropane	50.00	51.67	103	71-130
1,3,5-Trimethylbenzene	50.00	47.97	96	76-122
1,3-Dichlorobenzene	50.00	53.20	106	79-120
1,3-Dichloropropane	50.00	46.92	94	82-118
1,4-Dichlorobenzene	50.00	52.61	105	77-116
1,4-Dioxane	50.00	42.26	84	70-119
2,2-Dichloropropane	50.00	47.82	96	71-125
2-Butanone (MEK)	50.00	45.77	92	70-128
4-Chlorotoluene	50.00	52.93	106	77-117
4-Isopropyltoluene	50.00	40.84	82	71-124
4-Methyl-2-pentanone	50.00	42.39	85	75-121
Acetone	50.00	50.93	102	70-118
Benzene	50.00	50.47	101	77-118
Bromobenzene	50.00	53.48	107	79-116
Bromochloromethane	50.00	47.96	96	79-122

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/06/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1106 Location: Franklin, MA

Lab File ID: J32674B.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Bromodichloromethane	50.00	48.45	97	82-127
Bromoform	50.00	52.24	104	78-122
Bromomethane	50.00	51.33	103	70-130
Carbon disulfide	50.00	53.49	107	70-130
Carbon tetrachloride	50.00	52.22	104	77-125
Chlorobenzene	50.00	50.38	101	80-118
Chloroform	50.00	46.57	93	80-113
Chloromethane	50.00	50.11	100	70-130
cis-1,2-Dichloroethene	50.00	49.09	98	85-120
cis-1,3-Dichloropropene	50.00	45.86	92	79-116
Di-isopropyl ether (D	50.00	48.88	98	78-121
Dibromochloromethane	50.00	44.98	90	79-122
Dibromomethane	50.00	49.73	99	78-120
Ethyl tertiary-butyl	50.00	46.67	93	81-122
Ethylbenzene	50.00	51.69	103	84-123
Hexachlorobutadiene	50.00	49.52	99	70-121
Isopropylbenzene	50.00	50.39	101	78-120
m,p-Xylenes	100.0	112.5		75-129
Methyl tert-butyl eth	50.00	44.74	89	70-127
Methylene chloride	50.00	46.85	94	72-128
n-Butylbenzene	50.00	44.20	88	70-124
n-Propylbenzene	50.00	49.80	100	80-127
Naphthalene	50.00	49.70	99	70-126
o-Xylene	50.00	50.68	101	78-118
sec-Butylbenzene	50.00	44.36	89	78-118
Styrene	50.00	51.55	103	83-124
tert-Butylbenzene	50.00	45.07	90	76-118
Tertiary-amyl methyl	50.00	47.56	95	81-123

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/06/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1106 Location: Franklin, MA

Lab File ID: J32674B.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Tetrachloroethene (PCE)	50.00	47.10	94	77-122
Toluene	50.00	48.05	96	78-120
trans-1,2-Dichloroethene	50.00	49.05	98	80-120
trans-1,3-Dichloropro	50.00	45.69	91	71-111
Trichloroethene (TCE)	50.00	49.95	100	74-119
Vinyl chloride	50.00	47.84	96	70-116

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 4 8260B Method Blank Summary

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Lab File ID: J32677.D Lab Sample ID: VBLK1106

Matrix: Water Date Analyzed: 11/06/07

Instrument ID: MS8 Date Extracted:

Time Analyzed: 1111

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711143-1	841071101-03	J32685.D	11/06/2007
2	E711143-2A	841071101-04	J32680.D	11/06/2007
3	VLCS1106	VLCS1106	J32674B.D	11/06/2007
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AD9E750

FORM 3 Water MADEP VPH Lab Control Sample

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/06/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1106 Location: Franklin, MA

Lab File ID: 2110602.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Benzene	25.00	25.55	102	70-130
Ethylbenzene	25.00	25.38	102	70-130
m,p-Xylenes	50.00	50.30	101	70-130
Methyl tert-butyl eth	25.00	24.64	98	70-130
Naphthalene	25.00	23.29	93	70-130
o-Xylene	25.00	24.75	99	70-130
Toluene	25.00	25.14	100	70-130
Benzene #2	0.0	23.28	0.0	_
Ethylbenzene #2	0.0	26.06	0.0	-
Naphthalene #2	0.0	26.03	0.0	-
Toluene #2	0.0	24.56	0.0	_

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Water MADEP VPH Lab Control Sample

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/07/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: VLCS1106 DUP Location: Franklin, MA

Lab File ID: 2110610.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec#	Rec
Benzene	25.00	25.42	102	70-130
Ethylbenzene	25.00	25.34	101	70-130
m,p-Xylenes	50.00	50.32	101	70-130
Methyl tert-butyl eth	25.00	24.96	100	70-130
Naphthalene	25.00	25.32	101	70-130
o-Xylene	25.00	24.43	98	70-130
Toluene	25.00	26.27	105	70-130
Benzene #2	0.0	26.90	0.0	_
Ethylbenzene #2	0.0	27.68	0.0	_
Naphthalene #2	0.0	29.47	0.0	_
Toluene #2	0.0	29.29	0.0	_

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 4 MADEP VPH Method Blank Summary

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Lab File ID: 2110604.D Lab Sample ID: VBLK1106

Matrix: Water Date Analyzed: 11/06/07

Instrument ID: GC2 Date Extracted:

Time Analyzed: 1611

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711143-2B	841071101-04	2110605.D	11/06/2007
2	E711143-3B	841071101-05	2110606.D	11/06/2007
3	E711143-4B	841071101-06	2110607.D	11/06/2007
4	VLCS1106	VLCS1106	2110602.D	11/06/2007
5	VLCS1106 DUP	VLCS1106 DUP	2110610.D	11/07/2007
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FORM 3 Soil MADEP EPH Lab Control Sample

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/07/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1106S-1 Location: Franklin, MA

Lab File ID: A23891.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	()	()	Rec#	Rec
2-Methylnaphthalene	4000	2027	51	40-140
Acenaphthene	4000	2152	54	40-140
Acenaphthylene	4000	2274	57	40-140
Anthracene	4000	2874	72	40-140
Benzo[a]anthracene	4000	3242	81	40-140
Benzo[a]pyrene	4000	3184	80	40-140
Benzo[b]fluoranthene	4000	3124	78	40-140
<pre>Benzo[g,h,i]perylene</pre>	4000	3179	79	40-140
Benzo[k]fluoranthene	4000	3201	80	40-140
Chrysene	4000	3064	77	40-140
Dibenz[a,h]anthracene	4000	3173	79	40-140
Fluoranthene	4000	2750	69	40-140
Fluorene	4000	2312	58	40-140
<pre>Indeno[1,2,3-cd]pyrene</pre>	4000	3128	78	40-140
Naphthalene	4000	1911	48	40-140
Phenanthrene	4000	2604	65	40-140
Pyrene	4000	2733	68	40-140

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Soil MADEP EPH Lab Control Sample Duplicate

Lab Name: Premier Laboratory, Inc Date Analyzed: 11/07/07

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Sample No.: LCS1106S-1 Location: Franklin, MA

Lab File ID: A23892.D

	Spike	Sample				QC
	Added	Concentration	%		Li	mits
Compound	()	()	Rec#	RPD#	RPD	Rec
2-Methylnaphthalene	4000	1662	42	19.4	25	40-140
Acenaphthene	4000	1736	43	22.7	25	40-140
Acenaphthylene	4000	1811	45	23.5	25	40-140
Anthracene	4000	2312	58	21.5	25	40-140
Benzo[a]anthracene	4000	2435	61	28.2*	25	40-140
Benzo[a]pyrene	4000	2433	61	27.0*	25	40-140
Benzo[b]fluoranthene	4000	2360	59	27.7*	25	40-140
<pre>Benzo[g,h,i]perylene</pre>	4000	2345	59	29.0*	25	40-140
Benzo[k]fluoranthene	4000	2280	57	33.6*	25	40-140
Chrysene	4000	2348	59	26.5*	25	40-140
Dibenz[a,h]anthracene	4000	2316	58	30.6*	25	40-140
Fluoranthene	4000	2111	53	26.2*	25	40-140
Fluorene	4000	1860	46	23.1	25	40-140
<pre>Indeno[1,2,3-cd]pyrene</pre>	4000	2289	57	31.1*	25	40-140
Naphthalene	4000	1617	40	18.2	25	40-140
Phenanthrene	4000	2009	50	26.1*	25	40-140
Pyrene	4000	2137	53	24.8	25	40-140

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 4 MADEP EPH Method Blank Summary

Project No.: E711143 Project: 20050458.B10/Nu-Style Phase

Lab File ID: A23893.D Lab Sample ID: E1106BS-1

Matrix: Soil Date Analyzed: 11/07/07

Instrument ID: GC1 Date Extracted:

Time Analyzed: 1923

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	LCS1106S-1	LCS1106S-1	A23891.D	11/07/2007
2	LCSD1106S-1	LCSD1106S-1	A23892.D	11/07/2007
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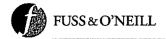
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☐ Other _

СНА	IN-OF-CUSTO	DY RECO	RD	14719		□ 1 Day* □ 3 Da □ 2 Days* → Stan	Turnacound iys* ====================================	1 Other (days) *Surcharge Applies
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Modified Tier II Data Validation Narrative and Certification

Project: 20050458B10, Former Nu-Style Company, Inc. Facility

Premier Laboratory Project Number:	E711585
Date Samples Received at Laboratory:	11/8/2007
Date of Review:	12/14/2007

Ten groundwater samples, including two field duplicates, were collected from eight monitoring wells over two days. Five of the wells were sampled using low-flow methodology; the remaining five had poor recharge and therefore grab samples were collected. Samples were submitted to Premier Laboratory in Dayville, Connecticut for analysis of RCRA-8 metals by EPA Methods 6010B and 245.2, volatile organic compounds (VOCs) by EPA Method 8260B, and/or petroleum hydrocarbons by Massachusetts Extractable and Volatile Petroleum Hydrocarbon (EPH/VPH) methods of analysis. Analyses for individual samples were determined based upon constituents of concern in the area where the individual monitoring wells are located.

Dedicated sampling equipment was used; therefore, no equipment blank was indicated. Two aqueous trip blanks were submitted, one for each day of sampling. No VOCs were reported in either trip blank.

Results of primary and duplicate sample pairs were generally similar. The relative percent difference (RPD) calculated for lead in one sample pair was 46% which is above the 30% limit established by the QAPP for aqueous samples.

Surrogate recoveries were acceptable for all applicable analyses. With the exception of hexachlorobutadiene, reporting limits were low enough to compare to MADEP GW-2 and GW-3 criteria. The lowest achievable reporting limit for this compound is $0.5~\mu g/L$.

I certify that the field and laboratory data associated with the above referenced project, to the best of my knowledge with the exceptions noted above, are compliant with the Quality Assurance Project Plan for the Former Nu-Style Company, Inc. Facility located in Franklin, Massachusetts dated September 2006.

Certified by:

Lynne P. Matteson QA/QC Officer



INITIAL DATE: MAY 2006 REVISION DATE: MAY 2006 REVISION: 0.0

PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY MODIFIED TIER I COMPLETENESS CHECKLIST

	<u>YES</u>	<u>NO</u>
1. SAMPLING AND FIELD MEASUREMENTS:	•	
Field measurement calibration records		
Groundwater field measurements (if applicable)		
Soil sampling field measurements (if applicable)		□ N/A
Sediment sampling field measurements (if applicable)		□ NA
Surface water sampling field measurements (if applicable)		□ h(V-
Low-flow sampling field measurements (if applicable)		
Documentation of field activities	V	
Sample numbering and labeling		
Chain-of-Custody records	☑	
Trip blanks	⊻′,	
Duplicate samples		□
Equipment blanks		
Split samples (if any)		□ νμ-
2. LABORATORY MEASUREMENTS: Trip blanks Instrument blanks Laboratory control samples Duplicates samples Equipment blanks Matrix spike/matrix spike duplicates Analysis type Chain-of-Custody records Surrogate recoveries Sample Project Narratives Split samples (if any) TOTAL:		0
PERCEI	NT COMPLETE: _	100 %



61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

ANALYTICAL DATA REPORT

Report Number: E711585 Project: 20050458.B10/Nu-Style Phase II

prepared for:

Fuss & O'Neill 275 Promenade Street Providence, RI 02908

Attn: David Foss

Received Date: 11/8/2007 Report Date: 11/20/2007

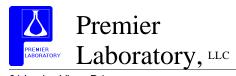
Premier Laboratory, LLC Authorized Signature



ADBA150

Certifications: CT (PH-0465), MA (M-CT008), ME (CT050), NH (2020), NJ (CT002), NY (11549), RI (RI246)

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		MADEP MCP A	Analytical Method Re	port Certifica	tion Form				
Labo	ratory Name: Prer	nier Laboratory, LL0	C		Project #:	E7115	35		
Proje	ct Location: Frank	klin, MA			MADEP RT	N ¹ :			
This	Form provides certif	fications for the follo	owing data set:[list La	boratory San	nple ID Numbe	er(s)]			
1, 10	, 11, 12, 2, 3, 4, 5,	6, 7, 8, 9							
		Groundwater □ S	oil/Sediment Drir	nking Water	☐ Other				
MCP	SW-846	8260B ⊠	8151A □	8330	□ 6010	OB ⊠			0A/1A □
Meth	ods Used	8270C □	8081A □	VPH		20 🗆		90	14M ² □
	ecified in MADEP	8082 🗆	8021B □	EPH	⊠ 7000 \$	S ³ □		7	7196A □
	endium of ical Methods.		g Number (RTN), if know 9014 or MADEP Physiolo		Cvanide (PAC) N	/lethod			
	all that apply)		7000 Series List individu			viculou			
	An affirmative res	ponse to question	s A, B, C, and D is I	required for	"Presumptive	Certail	nty"	sta	tus
Α			oratory in a condition			X Y			No ¹
	1	•	dy documentation for						
В			for the specified ana			⊠ Y	es		No ¹
	included in this rep	oort followed, includ	ling the requirement t	o note and					
	discuss in a narrat	tive QC data that di	d not meet appropria	te performan	ce				
	standards or guide	elines?							
С	Does the analytica	al data included in t	his report meet all the	e requiremen	nts	⊠ Y	es		No ¹
	for "Presumptive C	Certainty", as descri	bed in Section 2.0 (a),(b),(c) and	(d) of the				
	MADEP documen	t CAM VII A, "Quali	ity Assurance and Qu	uality Control	Guidelines				
	for the Acquisition	and Reporting of A	Analytical Data"?						
D	VPH and EPH Me	ethods only: Was t	the VPH or EPH met	hod run witho	out	⊠ Y	es		No ¹
	significant modific	ations, as specified	in Section 11.3?						
	A response	to questions E an	nd F below is require	ed for "Presu	umptive Certa	inty" st	atus	s	
Е	Were all QC perfo	rmance standards a	and recommendations	s for the		⊠ Y	es		No ¹
	specified methods	achieved?							
F	Were results for a	II analyte-list compo	ounds/elements for th	ne specified		⊠ Y	es		No ¹
	method(s) reported	d?							
¹ All NO answers must be addressed in an attached Environmental Laboratory case narrative.									
·									
I, the undersigned, attest under the pains and penalties of perjury that, based upon my personal inquiry of those responsible for obtaining the information, the material contained in this									
analytical report is, to the best of my knowledge and belief, accurate and complete.									
Signature: Position: Laboratory Director									
Print	ed Name: Robert	t Stevenson		Date: 11/2	20/2007				
II .									

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61 Louisa Viens Drive Dayville, CT 06241 FAX: 860-774-2689 860-774-6814 800-932-1150

Report No: E711585 Client: Fuss & O'Neill

Project: 20050458.B10/Nu-Style Phase II

CASE NARRATIVE / METHOD CONFORMANCE SUMMARY

Premier Laboratory received 12 samples from Fuss & O'Neill on 11/08/2007. The samples were analyzed from the following list of analyses:

Extractable Petroleum Hydrocarbon (EPH) MADEP EPH[MADEP EPH] Trace Metals by 6010B 6010B[3000] Volatiles by 8260B (GA/GW-1/S-1) 8260B Mercury by 245.2 in DW/WW 245.2[245.1] Volatile Petroleum Hydrocarbon (VPH) MADEP VPH

Variances:

SDG:

The lowest requested standard for Hexachlorobutadiene of 0.45 ug/L was not achieved. The lowest possible detection limit of 0.5 ug/L for Hexachlorobutadiene was reported.

Method:

None reported.

QA/QC:

None reported.

Laboratory: Premier Laboratory, LLC
PL Report No: E711585
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 11/8/2007 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(1) 841071106-01					
Date Collected: 11/6/2007 Matrix: Aque	POLIC				
Trace Metals by 6010B	zous .				
Arsenic	ND	0.0050	mg/L	11/15/07 13	3:51 AMM
Barium	0.060	0.0020	mg/L		3:51 AMM
Cadmium	ND	0.0020	mg/L		3:51 AMM
Chromium	ND	0.0020	mg/L		3:51 AMM
Lead	ND	0.0020	mg/L		3:51 AMM
Selenium	ND	0.0050	mg/L		3:51 AMM
Silver	ND	0.0020	mg/L		3:51 AMM
Mercury by 245.2 in DW/WW	ND	0.00020	mg/L	11/15/07	KAW
(2) 841071106-02					
Date Collected: 11/6/2007 Matrix: Aque	eous				
Trace Metals by 6010B					
Arsenic	ND	0.0050	mg/L	11/15/07 13	3:53 AMM
Barium	0.061	0.0020	mg/L	11/15/07 13	3:53 AMM
Cadmium	ND	0.0020	mg/L	11/15/07 13	3:53 AMM
Chromium	ND	0.0020	mg/L	11/15/07 13	3:53 AMM
Lead	ND	0.0020	mg/L	11/15/07 13	3:53 AMM
Selenium	ND	0.0050	mg/L	11/15/07 13	3:53 AMM
Silver	ND	0.0020	mg/L	11/15/07 13	3:53 AMM
Mercury by 245.2 in DW/WW	ND	0.00020	mg/L	11/15/07	KAW
(3) 841071106-03					
Date Collected: 11/6/2007 Matrix: Aque	<u>eous</u>				
Trace Metals by 6010B					
Arsenic	ND	0.0050	mg/L		3:54 AMM
Barium	0.031	0.0020	mg/L	11/15/07 13	3:54 AMM
Cadmium	ND	0.0020	mg/L	11/15/07 13	3:54 AMM
Chromium	0.0029	0.0020	mg/L	11/15/07 13	3:54 AMM
Lead	0.0066	0.0020	mg/L		3:54 AMM
Selenium	ND	0.0050	mg/L		3:54 AMM
Silver	ND	0.0020	mg/L		3:54 AMM
Mercury by 245.2 in DW/WW	ND	0.00020	mg/L	11/15/07	KAW

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Laboratory: Premier Laboratory, LLC
PL Report No: E711585
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 11/8/2007 Project: 20050458.B10/Nu-Style Phase II

Matrix M	Parameter		Result	DL	Units	Completed		Ву Г	Dilution
Date Collected: 11/6/2007	(4) 941071104 04								
Trace Metals by 6010B		Matrix: Aqueous							
Arsenic		Matrix. Aqueous							
Barium	•		ND	0.0050	ma/I	11/15/07	13.56	ΛММ	
Cadmium					-				
Chromium					-				
Lead					-				
Selenium					-				
Silver					-				
Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 KAW					-				
Sample S					-		13.30		
Date Collected: 11/6/2007 Matrix: Aqueous	Mercury by 243.2 III DW/W W		ND	0.00020	mg/L	11/13/07		KAW	
Date Collected: 11/6/2007 Matrix: Aqueous	(5) 841071106 ₋ 05								
Trace Metals by 6010B		Matriv: Aqueous							
Arsenic ND 0.0050 mg/L 11/15/07 13:58 AMM Barium 0.24 0.0020 mg/L 11/15/07 13:58 AMM Cadmium ND 0.0020 mg/L 11/15/07 13:58 AMM Chromium ND 0.0020 mg/L 11/15/07 13:58 AMM Lead 0.0033 0.0020 mg/L 11/15/07 13:58 AMM Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0050 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 14:00 AMM Arsenic ND 0.0020 mg/L 11/15/07 14:00 AMM <tr< td=""><td></td><td>Matrix. Aqueous</td><td></td><td></td><td></td><td></td><td></td><td></td><td></td></tr<>		Matrix. Aqueous							
Barium 0.24 0.0020 mg/L 11/15/07 13:58 AMM Cadmium ND 0.0020 mg/L 11/15/07 13:58 AMM Chromium ND 0.0020 mg/L 11/15/07 13:58 AMM Lead 0.0033 0.0020 mg/L 11/15/07 13:58 AMM Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0050 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 14:00 AMM Mercury by 245.2 in DW/WW Matrix: Aqueous ND 11/15/07	•		ND	0.0050	ma/I	11/15/07	13.58	ΛММ	
Cadmium ND 0.0020 mg/L 11/15/07 13:58 AMM Chromium ND 0.0020 mg/L 11/15/07 13:58 AMM Lead 0.0033 0.0020 mg/L 11/15/07 13:58 AMM Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 14:00 XAW (6) 841071106-06 Date Collected: 11/6/2007 Matrix: Aqueous NB 0.0050 mg/L 11/15/07 14:00 AMM Arsenic ND <td></td> <td></td> <td></td> <td></td> <td>-</td> <td></td> <td></td> <td></td> <td></td>					-				
Chromium ND 0.0020 mg/L 11/15/07 13:58 AMM Lead 0.0033 0.0020 mg/L 11/15/07 13:58 AMM Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM Mean Call Matrix: Aqueous NB 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Lead 0.0033 0.0020 mg/L 11/15/07 13:58 AMM Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM MEAUTION STATES AQUEOUS ND 0.00020 mg/L 11/15/07 14:00 AMM Barium ND 0.0050 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00					-				
Selenium ND 0.0050 mg/L 11/15/07 13:58 AMM Silver ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.0020 mg/L 11/15/07 KAW (6) 841071106-06 Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B Arsenic ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Silver ND 0.0020 mg/L 11/15/07 13:58 AMM Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 13:58 AMM (6) 841071106-06 Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Mercury by 245.2 in DW/WW ND 0.00020 mg/L 11/15/07 KAW (6) 841071106-06 Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
(6) 841071106-06 Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B Arsenic ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-		13.30		
Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B Arsenic ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM	Mercury by 243.2 III DW/W W		ND	0.00020	mg/L	11/13/07		KAW	
Date Collected: 11/6/2007 Matrix: Aqueous Trace Metals by 6010B Arsenic ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM	(6) 841071106-06								
Trace Metals by 6010B Arsenic ND 0.0050 mg/L 11/15/07 14:00 AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM		Matrix: Aqueous							
Arsenic ND 0.0050 mg/L 11/15/07 14:00 MM AMM Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM AMM		1,2442 11940 045							
Barium 0.25 0.0020 mg/L 11/15/07 14:00 AMM AMM Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM	•		ND	0.0050	mg/L	11/15/07	14:00	AMM	
Cadmium ND 0.0020 mg/L 11/15/07 14:00 AMM Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Chromium ND 0.0020 mg/L 11/15/07 14:00 AMM Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Lead 0.0053 0.0020 mg/L 11/15/07 14:00 AMM Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Selenium ND 0.0050 mg/L 11/15/07 14:00 AMM Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
Silver ND 0.0020 mg/L 11/15/07 14:00 AMM					-				
č					-				
	Mercury by 245.2 in DW/WW		ND	0.00020	mg/L	11/15/07		KAW	

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Laboratory: Premier Laboratory, LLC
PL Report No: E711585
Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 11/8/2007 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(8) 841071107-08					
Date Collected: 11/7/2007 Matrix: Aqueous					
Trace Metals by 6010B					
Arsenic	ND	0.0050	mg/L	11/15/07 14:20	AMM
Barium	0.11	0.0020	mg/L	11/15/07 14:20	AMM
Cadmium	ND	0.0020	mg/L	11/15/07 14:20	AMM
Chromium	ND	0.0020	mg/L	11/15/07 14:20	AMM
Lead	ND	0.0020	mg/L	11/15/07 14:20	AMM
Selenium	ND	0.0050	mg/L	11/15/07 14:20	AMM
Silver	ND	0.0020	mg/L	11/15/07 14:20	AMM
Mercury by 245.2 in DW/WW	ND	0.00020	mg/L	11/15/07	KAW
(9) 841071107-09					
Date Collected: 11/7/2007 Matrix: Aqueous					
Trace Metals by 6010B, Dissolved					
Arsenic	ND	0.0050	mg/L	11/15/07	AMM
Barium	0.39	0.0020	mg/L	11/15/07	AMM
Cadmium	ND	0.0020	mg/L	11/15/07	AMM
Chromium	ND	0.0020	mg/L	11/15/07	AMM
Lead	0.094	0.0020	mg/L	11/15/07	AMM
Selenium	ND	0.0050	mg/L	11/15/07	AMM
Silver	ND	0.0020	mg/L	11/15/07	AMM
Mercury by 245.2 in DW/WW, Dissolved	ND	0.00020	mg/L	11/15/07	KAW
(10) 841071107-10					
Date Collected: 11/7/2007 Matrix: Aqueous					
Trace Metals by 6010B, Dissolved					
Arsenic	ND	0.0050	mg/L	11/15/07	AMM
Barium	0.18	0.0020	mg/L	11/15/07	AMM
Cadmium	ND	0.0020	mg/L	11/15/07	AMM
Chromium	ND	0.0020	mg/L	11/15/07	AMM
Lead	0.0026	0.0020	mg/L	11/15/07	AMM
Selenium	ND	0.0050	mg/L	11/15/07	AMM
Silver	ND	0.0020	mg/L	11/15/07	AMM
Mercury by 245.2 in DW/WW, Dissolved	ND	0.00020	mg/L	11/15/07	KAW

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Laboratory: Premier Laboratory, LLC

PL Report No: E711585

Customer: Fuss & O'Neill
Location: Franklin, MA

Date Received: 11/8/2007 Project: 20050458.B10/Nu-Style Phase II

Parameter	Result	DL	Units	Completed	By Dilution
(11) 841071107-11 <u>Date Collected: 11/7/2007 Matrix: Aqueous</u>					
Trace Metals by 6010B, Dissolved	N.D.	0.0050	77	11/15/05	43.07
Arsenic	ND	0.0050	mg/L	11/15/07	AMM
Barium	0.17	0.0020	mg/L	11/15/07	AMM
Cadmium	ND	0.0020	mg/L	11/15/07	AMM
Chromium	ND	0.0020	mg/L	11/15/07	AMM
Lead	0.0060	0.0020	mg/L	11/15/07	AMM
Selenium	ND	0.0050	mg/L	11/15/07	AMM
Silver	ND	0.0020	mg/L	11/15/07	AMM
Mercury by 245.2 in DW/WW, Dissolved	ND	0.00020	mg/L	11/15/07	KAW

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841071106-01
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Analyzed:	-11/16/07	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene #2	93	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Matrix: Aqueous

Percent Moisture: N/A

Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E711585 PL Sample No: 1 Sample Description: 841071106-01

Date Collected: 11/6/2007 Date Received: 11/8/2007 Date Extracted: By: Date Analyzed: 11/13/07 By: DDD

Method: 8260B QC Batch#: 57819

Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32815.D Units: ug/L

167-64-1	CAS No.	Parameter	Result	DL
Total	67-64-1	Acetone	ND	10
108-86-1 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 0.50 75-27-4 Bromochloromethane ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 ND 1.0 75-11-8 nButylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-62-3-5 Carbon tetrachloride ND 1.0 75-60-3 Chlorotehane ND 1.0 75-00-3 Chlorotehane ND 1.0 75-00-3 Chlorotehane ND 1.0 74-87-3 Chlorotehane ND 1.0 74-87-3 Chlorotehane ND 1.0 74-87-3 Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 7	107-13-1	Acrylonitrile	ND	0.50
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 0.50 75-27-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n.Butylbenzene ND 1.0 104-51-8 n.Butylbenzene ND 1.0 98-06-6 tetr-Butylbenzene ND 1.0 75-51-50 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-40-3 Chlorobenzene ND 1.0 74-87-3 Chlorotofuene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 96-12-8 1,2-Dichlorobenzene ND 0.50 <t< td=""><td>71-43-2</td><td>Benzene</td><td>ND</td><td>1.0</td></t<>	71-43-2	Benzene	ND	1.0
75-27-4 Bromofchorom ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorofene ND 1.0 108-90-7 Chlorofene ND 1.0 75-0-3 Chloroferm ND 1.0 74-87-3 Chlorofene ND 1.0 74-87-3 Chlorofene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.5 106-93-4 4-Chlorotoluene ND 0.5 106-93-4 1,2-Dibromo-directhane ND 0.5 106-93-4 1,2-Dibromo-directhane ND 0.5 106-93-4 </td <td>108-86-1</td> <td>Bromobenzene</td> <td>ND</td> <td>1.0</td>	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chloroform ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 14-95-3 Dibromochloromethane ND 0.50	74-97-5	Bromochloromethane	ND	1.0
Part	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloroteme ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 1.0 95-50-1 1,2-Dibromochloromethane ND 1.0 94-95-3 Dibromochloromethane ND 1.0 <	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 1.0 98-06-6 tert-Butylbenzene ND 1.0	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon terrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorothane ND 1.0 67-66-3 Chlorofform ND 1.0 74-87-3 Chlorothane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND <td< td=""><td>104-51-8</td><td></td><td>ND</td><td>1.0</td></td<>	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorofeme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 10-46-7 1,4-Dichlorobenzene ND <td>135-98-8</td> <td>•</td> <td>ND</td> <td>1.0</td>	135-98-8	•	ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroferm ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorocthane ND	98-06-6		ND	1.0
108-90-7				1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-schloromethane ND 0.50 149-5-3 Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroptopene ND	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 1495-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichlorotethane ND 1.0 107-06-2 1,2-Dichloroptethane ND 1.0 75-35-4 1,1-Dichloroptethene ND 1.0 156-69-2 cis-1,2-Dichloroptropane ND 1.0 18-87-5 1,2-Dichloroptropane <td< td=""><td>108-90-7</td><td>Chlorobenzene</td><td>ND</td><td>1.0</td></td<>	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-69-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane </td <td>75-00-3</td> <td>Chloroethane</td> <td>ND</td> <td>1.0</td>	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropo	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 107-06-2 1,2-Dichlorothane ND 1.0 75-35-4 1,1-Dichlorothene ND 1.0 156-59-2 cis-1,2-Dichlorothene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropro	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 540-64-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 <	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dic	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropr	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	124-48-1		ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9		ND	1.0
563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7		ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6		ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7		ND	1.0
	123-91-1	•	ND	20

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Matrix: Aqueous

Percent Moisture: N/A Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E711585 PL Sample No: 1 (continued) Sample Description: 841071106-01

Date Collected: 11/6/2007 Date Received: 11/8/2007 Date Extracted: By: Date Analyzed: 11/13/07 By: DDD

Dilution Factor: 1 Method: 8260B Soil Extract Volume: Lab Data File: J32815.D QC Batch#: 57819

Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene ND		1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene		1.0
75-01-4	Vinyl chloride N		1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	m,p-Xylenes ND	
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	101%	89%-113%	
Bromofluorobenzene	86%	83%-107%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	101%	89%-113%
Bromofluorobenzene	86%	83%-107%
Toluene-d8	99%	88%-108%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841071106-02
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Analyzed:	-11/16/07	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene #2	100	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes

^{***} Excludes Naphthalene

Matrix: Aqueous

Percent Moisture: N/A

Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 Sample Description: 841071106-02

Date Collected: 11/6/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/13/07 By: DDD

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57819
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32816.D

Units: ug/L

167-64-1	CAS No.	Parameter	Result	DL
Total	67-64-1	Acetone	ND	10
108-86-1 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 0.50 75-27-4 Bromochloromethane ND 0.50 75-27-4 Bromofichloromethane ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 ND 1.0 74-83-9 ND 1.0 75-11-8 NB NB NB NB 1.0 75-11-8 ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 74-87-3 Chlorothane ND 1.0 74-87-3 Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-1 75-71-8 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-73-4 1.1-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 1.0 75-73-5 1.2-Dichlorothane ND 1.0	107-13-1	Acrylonitrile	ND	0.50
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 0.50 75-27-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n.Butylbenzene ND 1.0 104-51-8 n.Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-51-50 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-40-3 Chlorobenzene ND 1.0 74-87-3 Chlorotofuene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 96-12-8 1,2-Dichlorobenzene ND 1.0 <tr< td=""><td>71-43-2</td><td>Benzene</td><td>ND</td><td>1.0</td></tr<>	71-43-2	Benzene	ND	1.0
75-27-4 Bromofchorom ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorofene ND 1.0 108-90-7 Chlorofene ND 1.0 75-0-3 Chloroferm ND 1.0 74-87-3 Chlorofene ND 1.0 74-87-3 Chlorofene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.5 106-93-4 4-Chlorotoluene ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 </td <td>108-86-1</td> <td>Bromobenzene</td> <td>ND</td> <td>1.0</td>	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chloroform ND 1.0 108-90-7 Chloroferme ND 1.0 75-00-3 Chloroform ND 1.0 74-87-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-	74-97-5	Bromochloromethane	ND	1.0
Part	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloroteme ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 1.0 95-50-1 1,2-Dibromochloromethane ND 1.0 94-95-3 Dibromochloromethane ND 1.0 <	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 1.0 98-06-6 tert-Butylbenzene ND 1.0	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon terrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorothane ND 1.0 67-66-3 Chlorofform ND 1.0 74-87-3 Chlorothane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND <td< td=""><td>104-51-8</td><td></td><td>ND</td><td>1.0</td></td<>	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorofeme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 10-46-7 1,4-Dichlorobenzene ND <td>135-98-8</td> <td>•</td> <td>ND</td> <td>1.0</td>	135-98-8	•	ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroferm ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorocthane ND	98-06-6		ND	1.0
108-90-7				1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-schloromethane ND 0.50 149-5-3 Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroptopene ND	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 1495-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichlorotethane ND 1.0 107-06-2 1,2-Dichloroptethane ND 1.0 75-35-4 1,1-Dichloroptethene ND 1.0 156-69-2 cis-1,2-Dichloroptropane ND 1.0 18-87-5 1,2-Dichloroptropane <td< td=""><td>108-90-7</td><td>Chlorobenzene</td><td>ND</td><td>1.0</td></td<>	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-69-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane </td <td>75-00-3</td> <td>Chloroethane</td> <td>ND</td> <td>1.0</td>	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropo	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 107-06-2 1,2-Dichlorothane ND 1.0 75-35-4 1,1-Dichlorothene ND 1.0 156-59-2 cis-1,2-Dichlorothene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropro	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 540-64-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 <	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dic	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropr	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	124-48-1		ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9		ND	1.0
563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7		ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6		ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7		ND	1.0
	123-91-1	•	ND	20

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 2 (continued) Sample Description: 841071106-02

Date Collected:11/6/2007Matrix: AqueousDate Received:11/8/2007Percent Moisture: N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By: DDDDilution Factor: 1

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1

Method: 8260B
QC Batch#: 57819
Units: ug/L

Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32816.D

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)		1.0
75-69-4	Trichlorofluoromethane ND		1.0
95-63-6	1,2,4-Trimethylbenzene ND		1.0
108-67-8	1,3,5-Trimethylbenzene ND 1.		1.0
75-01-4	Vinyl chloride ND 1.		1.0
95-47-6	· · · · · · · · · · · · · · · · · · ·		1.0
108-38-3	· · · · · · · · · · · · · · · · · · ·		1.0
Surrogate	Recovery Lin	nits	
1.0 D' 11 4 14	1020/ 000	1120/	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	103%	89%-113%
Bromofluorobenzene	87%	83%-107%
Toluene-d8	93%	88%-108%

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VOLATILE PETROLEUM HYDROCARBON (VPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841071106-03
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Analyzed:	-11/16/07	Percent Moisture:	N/A
		Method:	MADEP VPH
		Ext Method:	5030B

(VPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C5-C8 Aliphatics*	1	ND	100	ug/L
C9-C12 Aliphatics**	1	ND	100	ug/L
C9-C10 Aromatics***	1	ND	100	ug/L

^{*} Excludes MTBE, Benzene, and Toluene

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
2,5-dibromotoluene #2	93	70%-130%

TARGETED VPH ANALYTES

Analyte	Results	$\mathbf{Q}\mathbf{L}$	Units
Benzene	ND	5.0	ug/L
Ethylbenzene	ND	5.0	ug/L
Methyl tert-butyl ether (MTBE)	ND	1.0	ug/L
Naphthalene	ND	5.0	ug/L
Toluene	ND	5.0	ug/L
m,p-Xylenes	ND	5.0	ug/L
o-Xylene	ND	5.0	ug/L

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^{**} Excludes Ethylbenzene, Xylenes *** Excludes Naphthalene

Matrix: Aqueous Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 Sample Description: 841071106-03

Date Collected: 11/6/2007

Date Received: 11/8/2007

Date Extracted: By:
Date Analyzed: 11/13/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By:DDDDilution Factor:1Method:8260BSoil Extract Volume:QC Batch#:57819Lab Data File:J32817.D

Units: ug/L

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 3 (continued) Sample Description: 841071106-03

Date Collected: 11/6/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/13/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57819Lab Data File:J32817.D

Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery I	Limits	
1,2-Dichloroethane-d4	105%	39%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	105%	89%-113%
Bromofluorobenzene	86%	83%-107%
Toluene-d8	98%	88%-108%

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Matrix: Aqueous

Percent Moisture: N/A

Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 Sample Description: 841071106-04

Date Collected: 11/6/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/13/07 By: DDD

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57819
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32818.D

Units: ug/L

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20
	,		

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 4 (continued) Sample Description: 841071106-04

Date Collected: 11/6/2007 Matrix: Aqueous

Date Received: 11/8/2007 Percent Moisture: N/A

Date Extracted: By: Sample Weight/Volume:

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1

Method: 8260P

Method: 8260B Soil Extract Volume: QC Batch#: 57819 Lab Data File: J32818.D

Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	12	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	20	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lin	nits	
1.2-Dichloroethane-d4	105% 899	%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	105%	89%-113%
Bromofluorobenzene	89%	83%-107%
Toluene-d8	96%	88%-108%

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Matrix: Aqueous

Percent Moisture: N/A Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 Sample Description: 841071106-05

Date Collected: 11/6/2007

Date Received: 11/8/2007

Date Extracted: By:
Date Analyzed: 11/13/07 By: DDD

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1

Method: 8260B

QC Batch#: 57819

Lab Data File: J32819.D

Units: ug/L

167-64-1	CAS No.	Parameter	Result	DL
Total	67-64-1	Acetone	ND	10
108-86-1 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 0.50 75-27-4 Bromochloromethane ND 0.50 75-27-4 Bromofichloromethane ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 ND 1.0 74-83-9 ND 1.0 75-11-8 NB NB NB NB 1.0 75-11-8 ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 74-87-3 Chlorothane ND 1.0 74-87-3 Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-1 75-71-8 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-73-4 1.1-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 1.0 75-73-5 1.2-Dichlorothane ND 1.0	107-13-1	Acrylonitrile	ND	0.50
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 0.50 75-27-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n.Butylbenzene ND 1.0 104-51-8 n.Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-51-50 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-40-3 Chlorobenzene ND 1.0 74-87-3 Chlorotofuene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 96-12-8 1,2-Dichlorobenzene ND 1.0 <tr< td=""><td>71-43-2</td><td>Benzene</td><td>ND</td><td>1.0</td></tr<>	71-43-2	Benzene	ND	1.0
75-27-4 Bromofchorom ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorofene ND 1.0 108-90-7 Chlorofene ND 1.0 75-0-3 Chloroferm ND 1.0 74-87-3 Chlorofene ND 1.0 74-87-3 Chlorofene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.5 106-93-4 4-Chlorotoluene ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 </td <td>108-86-1</td> <td>Bromobenzene</td> <td>ND</td> <td>1.0</td>	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chloroform ND 1.0 108-90-7 Chloroferme ND 1.0 75-00-3 Chloroform ND 1.0 74-87-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-	74-97-5	Bromochloromethane	ND	1.0
Part	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloroteme ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 1.0 95-50-1 1,2-Dibromochloromethane ND 1.0 94-95-3 Dibromochloromethane ND 1.0 <	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 1.0 98-06-6 tert-Butylbenzene ND 1.0	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon terrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorothane ND 1.0 67-66-3 Chlorofform ND 1.0 74-87-3 Chlorothane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND <td< td=""><td>104-51-8</td><td></td><td>ND</td><td>1.0</td></td<>	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorofeme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 10-46-7 1,4-Dichlorobenzene ND <td>135-98-8</td> <td>•</td> <td>ND</td> <td>1.0</td>	135-98-8	•	ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroferm ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorocthane ND	98-06-6		ND	1.0
108-90-7				1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-schloromethane ND 0.50 149-5-3 Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroptopene ND	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 1495-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichlorotethane ND 1.0 107-06-2 1,2-Dichloroptethane ND 1.0 75-35-4 1,1-Dichloroptethene ND 1.0 156-69-2 cis-1,2-Dichloroptropane ND 1.0 18-87-5 1,2-Dichloroptropane <td< td=""><td>108-90-7</td><td>Chlorobenzene</td><td>ND</td><td>1.0</td></td<>	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-69-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane </td <td>75-00-3</td> <td>Chloroethane</td> <td>ND</td> <td>1.0</td>	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropo	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 107-06-2 1,2-Dichlorothane ND 1.0 75-35-4 1,1-Dichlorothene ND 1.0 156-59-2 cis-1,2-Dichlorothene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropro	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 540-64-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 <	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dic	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropr	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	124-48-1		ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9		ND	1.0
563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7		ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6		ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7		ND	1.0
	123-91-1	•	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 5 (continued) Sample Description: 841071106-05

Date Collected: 11/6/2007
Date Received: 11/8/2007
Date Extracted: By:
Date Analyzed: 11/13/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57819Lab Data File:J32819.D

Units: ug/L

CAS No.	Parameter	Resu	lt DL
100-41-4	Ethylbenzene	N	D 1.0
87-68-3	Hexachlorobutadiene	N	D 0.50
591-78-6	2-Hexanone	N	D 5.0
98-82-8	Isopropylbenzene	N	D 1.0
99-87-6	4-Isopropyltoluene	N	D 1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	1	.5 1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	N	D 5.0
75-09-2	Methylene chloride	N	D 5.0
91-20-3	Naphthalene	N	D 1.0
103-65-1	n-Propylbenzene	N	D 1.0
100-42-5	Styrene	N	D 1.0
109-99-9	Tetrahydrofuran	N	D 1.0
110-57-6	trans-1,4-Dichloro-2-butene	N	D 5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	N	D 1.0
96-18-4	1,2,3-Trichloropropane	N	D 1.0
630-20-6	1,1,1,2-Tetrachloroethane	N	D 1.0
79-34-5	1,1,2,2-Tetrachloroethane	N	D 0.50
127-18-4	Tetrachloroethene (PCE)	29	00 1.0
108-88-3	Toluene	N	D 1.0
87-61-6	1,2,3-Trichlorobenzene	N	D 1.0
120-82-1	1,2,4-Trichlorobenzene	N	D 1.0
71-55-6	1,1,1-Trichloroethane	N	D 1.0
79-00-5	1,1,2-Trichloroethane	N	D 1.0
79-01-6	Trichloroethene (TCE)	(50 1.0
75-69-4	Trichlorofluoromethane	N	D 1.0
95-63-6	1,2,4-Trimethylbenzene	N	D 1.0
108-67-8	1,3,5-Trimethylbenzene	N	D 1.0
75-01-4	Vinyl chloride	N	D 1.0
95-47-6	o-Xylene	N	D 1.0
108-38-3	m,p-Xylenes	N	D 1.0
Surrogate	Recovery	Limits	
1,2-Dichloroethane-d4	102%	89%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	102%	89%-113%
Bromofluorobenzene	87%	83%-107%
Toluene-d8	98%	88%-108%

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Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E711585 PL Sample No: 6 Sample Description: 841071106-06

Date Collected: 11/6/2007 Date Received: 11/8/2007 Date Extracted: By: Date Analyzed: 11/14/07 By: DDD

Method: 8260B QC Batch#: 57868

Units: ug/L

Matrix: Aqueous Percent Moisture: N/A Sample Weight/Volume: Dilution Factor: 1 Soil Extract Volume: Lab Data File: J32827.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 6 (continued) Sample Description: 841071106-06

Date Collected: 11/6/2007
Date Received: 11/8/2007
Date Extracted: By:
Date Analyzed: 11/14/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/14/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57868Lab Data File:J32827.D

Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	1.4	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	260	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	56	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lir	nits	
1.2-Dichloroethane-d4	100% 899	%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	100%	89%-113%
Bromofluorobenzene	88%	83%-107%
Toluene-d8	99%	88%-108%

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 Sample Description: 841071106-07

Date Collected:11/6/2007Matrix: AqueousDate Received:11/8/2007Percent Moisture: N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By: DDDDilution Factor: 1

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57819
Units: ug/L
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32804.D

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 7 (continued) Sample Description: 841071106-07

Date Collected: 11/6/2007 Matrix: Aqueous
Date Received: 11/8/2007 Percent Moisture: N/A
Date Extracted: By: Sample Weight/Volume:
Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1

Date Analyzed: 11/13/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57819
Units: ug/L
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32804.D

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	ND	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lir	nits	
1.2 Dichloroothana d4	1020/ 900	0/. 1120/.	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	102%	89%-113%
Bromofluorobenzene	88%	83%-107%
Toluene-d8	96%	88%-108%

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 Sample Description: 841071107-08

Date Collected: 11/7/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/14/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/14/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57868Lab Data File:J32828.D

Units: ug/L

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 8 (continued) Sample Description: 841071107-08

Date Collected: 11/7/2007
Date Received: 11/8/2007
Date Extracted: By:
Date Analyzed: 11/14/07 By: DDD

Date Analyzed: 11/14/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57868
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32828.D

Units: ug/L

CAS No.	Parameter	R	esult	DL
100-41-4	Ethylbenzene		ND	1.0
87-68-3	Hexachlorobutadiene		ND	0.50
591-78-6	2-Hexanone		ND	5.0
98-82-8	Isopropylbenzene		ND	1.0
99-87-6	4-Isopropyltoluene		ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	5.0
75-09-2	Methylene chloride		ND	5.0
91-20-3	Naphthalene		ND	1.0
103-65-1	n-Propylbenzene		ND	1.0
100-42-5	Styrene		ND	1.0
109-99-9	Tetrahydrofuran		ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene		ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane		ND	1.0
96-18-4	1,2,3-Trichloropropane		ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane		ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane		ND	0.50
127-18-4	Tetrachloroethene (PCE)		41	1.0
108-88-3	Toluene		ND	1.0
87-61-6	1,2,3-Trichlorobenzene		ND	1.0
120-82-1	1,2,4-Trichlorobenzene		ND	1.0
71-55-6	1,1,1-Trichloroethane		ND	1.0
79-00-5	1,1,2-Trichloroethane		ND	1.0
79-01-6	Trichloroethene (TCE)		45	1.0
75-69-4	Trichlorofluoromethane		ND	1.0
95-63-6	1,2,4-Trimethylbenzene		ND	1.0
108-67-8	1,3,5-Trimethylbenzene		ND	1.0
75-01-4	Vinyl chloride		ND	1.0
95-47-6	o-Xylene		ND	1.0
108-38-3	m,p-Xylenes		ND	1.0
Surrogate	Recovery	Limits		
1,2-Dichloroethane-d4	103%	89%-113%		
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Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	103%	89%-113%
Bromofluorobenzene	88%	83%-107%
Toluene-d8	96%	88%-108%

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Matrix: Aqueous

Dilution Factor: 1

Soil Extract Volume:

Percent Moisture: N/A

Sample Weight/Volume:

Lab Data File: J32829.D

Laboratory: Premier Laboratory, LLC Customer: Fuss & O'Neill Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E711585 PL Sample No: 9 Sample Description: 841071107-09

Date Collected: 11/7/2007 Date Received: 11/8/2007 Date Extracted: By: Date Analyzed: 11/14/07 By: DDD

Method: 8260B QC Batch#: 57868

Units: ug/L

107.13-1	CAS No.	Parameter	Result	DL
Times	67-64-1	Acetone	ND	10
108-86-1 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 1.0 1.0 75-27-4 Bromochloromethane ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 1.0 135-98-8 sec-Butylbenzene ND 1.0	107-13-1	Acrylonitrile	ND	0.50
Table	71-43-2	Benzene	ND	1.0
75-27-4 Bromofichloromethane ND 0.50 75-25-2 Bromoform ND 1.0 75-25-2 Bromoform ND 1.0 75-25-2 Bromoferm ND 1.0 75-25-2 Bromoferm ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 105-98-8 sec-Butylbenzene ND 1.0 1.	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoferm ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 1	74-97-5	Bromochloromethane	ND	1.0
Part	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 55-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroform ND 1.0 76-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 14-95-3 Dibromochlane (EDB) ND 0.50 74-95-3 Dibromochlane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0	75-25-2	Bromoform	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 106-93-4 1,2-Dibromo-dhare (EDB) ND 0.50 106-93-4 1,2-Dibromethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 <td>74-83-9</td> <td>Bromomethane</td> <td>ND</td> <td>1.0</td>	74-83-9	Bromomethane	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 </td <td>78-93-3</td> <td>2-Butanone (MEK)</td> <td>ND</td> <td>5.0</td>	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorofore ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochlane (EDB) ND 0.50 106-93-4 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorofdiluoromethane ND 1.0	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 76-66-3 Chloroethane ND 1.0 74-87-3 Chloroethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-42-8 1,2-Dichlorobenzene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 106-43-4 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromo-chloromethane ND 0.50 124-48-1 Dibromo-chloromethane ND 0.50 124-49-3 Dibromo-thane (EDB) ND 0.50 14-495-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 175-31-8 Dichlorotifluoromethane ND 1.0 175-34-3 1,1-Dichloroethane ND 1.0 175-35-4 1,1-Dichloroethane ND 1.0 175-35-4 1,1-Dichloroethane ND 1.0 175-35-4 1,1-Dichloroethane ND 1.0 176-60-5 trans-1,2-Dichloroethene ND 1.0 186-60-5 trans-1,2-Dichloroethene ND 1.0 188-75- 1,2-Dichloroethene ND 1.0 188-75- 1,2-Dichloropopane ND 1.0 188-75- 1,2-Dichloropopane ND 1.0 190-20-7 2,2-Dichloropopane ND 1.0 1006-10-5 cis-1,3-Dichloropropane ND 1.0 1006-10-5 cis-1,3-Dichloropropane ND 1.0 1006-10-5 cis-1,3-Dichloropropane ND 0.50 10061-02-6 trans-1,3-Dichloropropane ND 0.50	135-98-8		ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroferm ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromoethane (EDB) ND 0.50 74-95-3 Dibromoethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorobenzene ND 1.0	98-06-6		ND	1.0
108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromoethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 75-35-4 1,1-Dichlorothane ND 1.0 75-35-4 1,1-Dichloropropane ND 1.	75-15-0		ND	1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 14-95-3 Dibromoethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane N	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroptopene ND 1.0 142-28-9 1,3-Dichloroptopane ND 1.0 590-20-7 2,2-Dichloroptopane	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromo-3-chloropropane (DBCP) ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 106-93-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-75 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloro	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropene ND 1.0 503-58-6 1,1-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dic	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroptopane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloroprop	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroptopane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloroprop	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether </td <td>124-48-1</td> <td>* * * * * * * * * * * * * * * * * * *</td> <td>ND</td> <td>0.50</td>	124-48-1	* * * * * * * * * * * * * * * * * * *	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether </td <td>106-93-4</td> <td>1,2-Dibromoethane (EDB)</td> <td>ND</td> <td>0.50</td>	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9	1,3-Dichloropropane	ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7	2,2-Dichloropropane	ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6		ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7	* *	ND	1.0
,	123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 9 (continued) Sample Description: 841071107-09

Date Collected:11/7/2007Matrix: AqueousDate Received:11/8/2007Percent Moisture: N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:11/14/07By: DDDDilution Factor: 1

Method: 8260B Soil Extract Volume: QC Batch#: 57868 Lab Data File: J32829.D

Units:	ug/L
Units:	ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	1.3	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	ND	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lin	nits	
1,2-Dichloroethane-d4	102% 899	%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	102%	89%-113%
Bromofluorobenzene	88%	83%-107%
Toluene-d8	100%	88%-108%

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Matrix: Aqueous

Percent Moisture: N/A

Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 10 Sample Description: 841071107-10

Date Collected: 11/7/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/14/07 By: DDD

Date Analyzed: 11/14/07 By: DDD Dilution Factor: 1
Method: 8260B
QC Batch#: 57868
Dilution Factor: 1
Soil Extract Volume:
Lab Data File: J32830.D

Units: ug/L

107-13-1 Acetone	CAS No.	Parameter	Result	DL
Total	67-64-1	Acetone	ND	10
108-86-1 Bromobenzene	107-13-1	Acrylonitrile	ND	0.50
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 0.50 75-27-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 104-51-8 n-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-40-3 Chloroberane ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 95-50-1 1,2-Dichlorobenzene ND 1.0	71-43-2	Benzene	ND	1.0
75-27-4 Bromofenm ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorofene ND 1.0 75-0-3 Chlorofene ND 1.0 75-6-3 Chlorofene ND 1.0 75-48-3 Chlorofene ND 1.0 74-87-3 Chlorofene ND 1.0 96-12-8 1.2-Dibromo-3-chloropropane (DBCP) ND 0.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1.2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-93-4 1.2-Dibromo-4-comethane ND 0.50 1	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tetrachloride ND 1.0 108-90-7 Chloroform ND 1.0 108-90-7 Chlorochtane ND 1.0 67-66-3 Chloroform ND 1.0 74-487-3 Chloroform ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 149-3 Dibromochloromethane ND 0.50	74-97-5	Bromochloromethane	ND	1.0
74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chloroethane ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 14-95-3 Dibromochlane (EDB) ND 0.50 74-95-3 Dibromochlane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 </td <td>75-27-4</td> <td>Bromodichloromethane</td> <td>ND</td> <td>0.50</td>	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 14-48-1 Dibromochloromethane ND 0.50 4-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 <td>75-25-2</td> <td>Bromoform</td> <td>ND</td> <td>1.0</td>	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chloroethane ND 1.0 75-00-3 Chloroform ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloroethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tertachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-3-chloropropane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene	135-98-8	· · · · · · · · · · · · · · · · · · ·	ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorotethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorothane ND 1.0 75-34-3 1,1-Dichlorothane ND <t< td=""><td>98-06-6</td><td></td><td>ND</td><td>1.0</td></t<>	98-06-6		ND	1.0
108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-brachen (EDB) ND 0.50 74-95-3 Dibromomethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloropropane ND </td <td></td> <td></td> <td></td> <td>1.0</td>				1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroptopane ND 1.0 156-60-5 trans-1,2-Dichloroptopane	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorodenane ND 1.0 107-06-2 1,2-Dichloropethane ND 1.0 75-35-4 1,1-Dichloropethene ND 1.0 156-69-2 cis-1,2-Dichloropethene ND 1.0 78-87-5 1,2-Dichloroperopane <td< td=""><td>108-90-7</td><td>Chlorobenzene</td><td>ND</td><td>1.0</td></td<>	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane <td>75-00-3</td> <td>Chloroethane</td> <td>ND</td> <td>1.0</td>	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromoethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dich	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 107-06-2 1,2-Dichlorothane ND 1.0 156-59-2 1,1-Dichlorothene ND 1.0 156-59-2 cis-1,2-Dichlorothene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dich	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 541-73-1 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-D	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroptopene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropro	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroptopane ND 1.0 78-87-5 1,2-Dichloroptopane ND 1.0 142-28-9 1,3-Dichloroptopane ND 1.0 563-58-6 1,1-Dichloroptopane ND 1.0 563-58-6 1,1-Dichloroptopane ND 0.50 10061-02-6 trans-1,3-Dichlorop	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether </td <td>124-48-1</td> <td></td> <td>ND</td> <td>0.50</td>	124-48-1		ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether </td <td>106-93-4</td> <td>1,2-Dibromoethane (EDB)</td> <td>ND</td> <td>0.50</td>	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene 1.7 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	1.7	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9		ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7	2,2-Dichloropropane	ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6	1,1-Dichloropropene	ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7	* *	ND	1.0
	123-91-1	1,4-Dioxane	ND	20

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Matrix: Aqueous

Percent Moisture: N/A Sample Weight/Volume:

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 10 (continued) Sample Description: 841071107-10

Date Collected: 11/7/2007
Date Received: 11/8/2007
Date Extracted: By:
Date Analyzed: 11/14/07 By: DDD

Date Analyzed: 11/14/07By: DDDDilution Factor: 1Method: 8260BSoil Extract Volume:QC Batch#: 57868Lab Data File: J32830.D

Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	23	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	25	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate	Recovery Lir	nits	
1.2-Dichloroethane-d4	100% 899	%-113%	

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	100%	89%-113%
Bromofluorobenzene	87%	83%-107%
Toluene-d8	96%	88%-108%

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Matrix: Aqueous

Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 11 Sample Description: 841071107-11

Date Collected: 11/7/2007

Date Received: 11/8/2007

Date Extracted: By:

Date Analyzed: 11/14/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/14/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57868Lab Data File:J32831.D

Units: ug/L

CAS No.	Parameter	Result	DL
67-64-1	Acetone	ND	10
107-13-1	Acrylonitrile	ND	0.50
71-43-2	Benzene	ND	1.0
108-86-1	Bromobenzene	ND	1.0
74-97-5	Bromochloromethane	ND	1.0
75-27-4	Bromodichloromethane	ND	0.50
75-25-2	Bromoform	ND	1.0
74-83-9	Bromomethane	ND	1.0
78-93-3	2-Butanone (MEK)	ND	5.0
104-51-8	n-Butylbenzene	ND	1.0
135-98-8	sec-Butylbenzene	ND	1.0
98-06-6	tert-Butylbenzene	ND	1.0
75-15-0	Carbon disulfide	ND	1.0
56-23-5	Carbon tetrachloride	ND	1.0
108-90-7	Chlorobenzene	ND	1.0
75-00-3	Chloroethane	ND	1.0
67-66-3	Chloroform	ND	1.0
74-87-3	Chloromethane	ND	1.0
95-49-8	2-Chlorotoluene	ND	1.0
106-43-4	4-Chlorotoluene	ND	1.0
96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
124-48-1	Dibromochloromethane	ND	0.50
106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
74-95-3	Dibromomethane	ND	1.0
95-50-1	1,2-Dichlorobenzene	ND	1.0
541-73-1	1,3-Dichlorobenzene	ND	1.0
106-46-7	1,4-Dichlorobenzene	ND	1.0
75-71-8	Dichlorodifluoromethane	ND	1.0
75-34-3	1,1-Dichloroethane	ND	1.0
107-06-2	1,2-Dichloroethane	ND	1.0
75-35-4	1,1-Dichloroethene	ND	1.0
156-59-2	cis-1,2-Dichloroethene	ND	1.0
156-60-5	trans-1,2-Dichloroethene	ND	1.0
78-87-5	1,2-Dichloropropane	ND	1.0
142-28-9	1,3-Dichloropropane	ND	1.0
590-20-7	2,2-Dichloropropane	ND	1.0
563-58-6	1,1-Dichloropropene	ND	1.0
10061-01-5	cis-1,3-Dichloropropene	ND	0.50
10061-02-6	trans-1,3-Dichloropropene	ND	0.50
60-29-7	Diethyl ether	ND	1.0
123-91-1	1,4-Dioxane	ND	20

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Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 11 (continued) Sample Description: 841071107-11

Date Collected:11/7/2007Matrix: AqueousDate Received:11/8/2007Percent Moisture: N/ADate Extracted:By:Sample Weight/Volume:Date Analyzed:11/14/07By: DDDDilution Factor: 1

Method: 8260B Soil Extract Volume:
QC Batch#: 57868 Lab Data File: J32831.D
Units: ug/L

CAS No.	Parameter	Result	DL
100-41-4	Ethylbenzene	ND	1.0
87-68-3	Hexachlorobutadiene	ND	0.50
591-78-6	2-Hexanone	ND	5.0
98-82-8	Isopropylbenzene	ND	1.0
99-87-6	4-Isopropyltoluene	ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)	ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)	ND	5.0
75-09-2	Methylene chloride	ND	5.0
91-20-3	Naphthalene	ND	1.0
103-65-1	n-Propylbenzene	ND	1.0
100-42-5	Styrene	ND	1.0
109-99-9	Tetrahydrofuran	ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene	ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane	ND	1.0
96-18-4	1,2,3-Trichloropropane	ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane	ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane	ND	0.50
127-18-4	Tetrachloroethene (PCE)	74	1.0
108-88-3	Toluene	ND	1.0
87-61-6	1,2,3-Trichlorobenzene	ND	1.0
120-82-1	1,2,4-Trichlorobenzene	ND	1.0
71-55-6	1,1,1-Trichloroethane	ND	1.0
79-00-5	1,1,2-Trichloroethane	ND	1.0
79-01-6	Trichloroethene (TCE)	59	1.0
75-69-4	Trichlorofluoromethane	ND	1.0
95-63-6	1,2,4-Trimethylbenzene	ND	1.0
108-67-8	1,3,5-Trimethylbenzene	ND	1.0
75-01-4	Vinyl chloride	ND	1.0
95-47-6	o-Xylene	ND	1.0
108-38-3	m,p-Xylenes	ND	1.0
Surrogate Recovery Limits			

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	102%	89%-113%
Bromofluorobenzene	88%	83%-107%
Toluene-d8	96%	88%-108%

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Matrix: Aqueous Percent Moisture: N/A

Laboratory: Premier Laboratory, LLC

Customer: Fuss & O'Neill
Location: Franklin, MA

PL Report No: E711585 Project: 20050458.B10/Nu-Style Phase II PL Sample No: 12 Sample Description: 841071107-12

Date Collected: 11/6/2007
Date Received: 11/8/2007
Date Extracted: By:
Date Analyzed: 11/13/07 By: DDD

Date Extracted:By:Sample Weight/Volume:Date Analyzed:11/13/07By:DDDMethod:8260BSoil Extract Volume:QC Batch#:57819Lab Data File:J32805.D

Units: ug/L

167-64-1	CAS No.	Parameter	Result	DL
Total	67-64-1	Acetone	ND	10
108-86-1 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 1.0 74-97-5 Bromochloromethane ND 0.50 75-27-4 Bromochloromethane ND 0.50 75-27-4 Bromofichloromethane ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 Bromomethane ND 1.0 74-83-9 ND 1.0 74-83-9 ND 1.0 75-11-8 NB NB NB NB 1.0 75-11-8 ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 75-00-3 Chlorothane ND 1.0 74-87-3 Chlorothane ND 1.0 74-87-3 Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-4 4-Chlorotoluene ND 1.0 74-88-1 75-71-8 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-71-8 Dichlorothenzene ND 1.0 75-73-4 1.1-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 75-73-5 1.2-Dichlorothane ND 1.0 1.0 75-73-5 1.2-Dichlorothane ND 1.0	107-13-1	Acrylonitrile	ND	0.50
74-97-5 Bromochloromethane ND 1.0 75-27-4 Bromodichloromethane ND 0.50 75-27-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n.Butylbenzene ND 1.0 104-51-8 n.Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-51-50 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-40-3 Chlorobenzene ND 1.0 74-87-3 Chlorotofuene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 106-43-4 4-Chlorotoluene ND 0.50 96-12-8 1,2-Dichlorobenzene ND 1.0 <tr< td=""><td>71-43-2</td><td>Benzene</td><td>ND</td><td>1.0</td></tr<>	71-43-2	Benzene	ND	1.0
75-27-4 Bromofchorom ND 0.50 75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chlorofene ND 1.0 108-90-7 Chlorofene ND 1.0 75-0-3 Chloroferm ND 1.0 74-87-3 Chlorofene ND 1.0 74-87-3 Chlorofene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.5 106-93-4 4-Chlorotoluene ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 1,2-Dibromo-directance ND 0.5 106-93-4 </td <td>108-86-1</td> <td>Bromobenzene</td> <td>ND</td> <td>1.0</td>	108-86-1	Bromobenzene	ND	1.0
75-25-2 Bromoform ND 1.0 74-83-9 Bromomethane ND 1.0 78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon disulfide ND 1.0 108-90-7 Chloroform ND 1.0 108-90-7 Chloroferme ND 1.0 75-00-3 Chloroform ND 1.0 74-87-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-	74-97-5	Bromochloromethane	ND	1.0
Part	75-27-4	Bromodichloromethane	ND	0.50
78-93-3 2-Butanone (MEK) ND 5.0 104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroteme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloroteme ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 1.0 95-50-1 1,2-Dibromochloromethane ND 1.0 94-95-3 Dibromochloromethane ND 1.0 <	75-25-2	Bromoform	ND	1.0
104-51-8	74-83-9	Bromomethane	ND	1.0
104-51-8 n-Butylbenzene ND 1.0 135-98-8 sec-Butylbenzene ND 1.0 1.0 98-06-6 tert-Butylbenzene ND 1.0	78-93-3	2-Butanone (MEK)	ND	5.0
135-98-8 sec-Butylbenzene ND 1.0 98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon terrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorothane ND 1.0 67-66-3 Chlorofform ND 1.0 74-87-3 Chlorothane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane (EDB) ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND <td< td=""><td>104-51-8</td><td></td><td>ND</td><td>1.0</td></td<>	104-51-8		ND	1.0
98-06-6 tert-Butylbenzene ND 1.0 75-15-0 Carbon disulfide ND 1.0 75-15-0 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chlorofeme ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chlorotoluene ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 124-48-1 Dibromochloromethane ND 0.50 74-95-3 Dibromochloromethane ND 0.50 74-95-3 Dibromochlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 10-46-7 1,4-Dichlorobenzene ND <td>135-98-8</td> <td>•</td> <td>ND</td> <td>1.0</td>	135-98-8	•	ND	1.0
75-15-0 Carbon disulfide ND 1.0 56-23-5 Carbon tetrachloride ND 1.0 108-90-7 Chlorobenzene ND 1.0 75-00-3 Chloroferm ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromochloromethane ND 0.50 74-95-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromoethane (EDB) ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorocthane ND	98-06-6		ND	1.0
108-90-7				1.0
75-00-3 Chloroethane ND 1.0 67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromo-schloromethane ND 0.50 149-5-3 Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroptopene ND	56-23-5	Carbon tetrachloride	ND	1.0
67-66-3 Chloroform ND 1.0 74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 1495-3 Dibromomethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichlorotethane ND 1.0 107-06-2 1,2-Dichloroptethane ND 1.0 75-35-4 1,1-Dichloroptethene ND 1.0 156-69-2 cis-1,2-Dichloroptropane ND 1.0 18-87-5 1,2-Dichloroptropane <td< td=""><td>108-90-7</td><td>Chlorobenzene</td><td>ND</td><td>1.0</td></td<>	108-90-7	Chlorobenzene	ND	1.0
74-87-3 Chloromethane ND 1.0 95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethene ND 1.0 156-69-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane </td <td>75-00-3</td> <td>Chloroethane</td> <td>ND</td> <td>1.0</td>	75-00-3	Chloroethane	ND	1.0
95-49-8 2-Chlorotoluene ND 1.0 106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropo	67-66-3	Chloroform	ND	1.0
106-43-4 4-Chlorotoluene ND 1.0 96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichlorothane ND 1.0 107-06-2 1,2-Dichlorothane ND 1.0 75-35-4 1,1-Dichlorothene ND 1.0 156-59-2 cis-1,2-Dichlorothene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropro	74-87-3	Chloromethane	ND	1.0
96-12-8 1,2-Dibromo-3-chloropropane (DBCP) ND 0.50 124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 540-64-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 142-28-9 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 <	95-49-8	2-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dic	106-43-4	4-Chlorotoluene	ND	1.0
124-48-1 Dibromochloromethane ND 0.50 106-93-4 1,2-Dibromoethane (EDB) ND 0.50 74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloropropane ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropane ND 1.0 503-58-6 1,1-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropr	96-12-8	1,2-Dibromo-3-chloropropane (DBCP)	ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	124-48-1		ND	0.50
74-95-3 Dibromomethane ND 1.0 95-50-1 1,2-Dichlorobenzene ND 1.0 541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 158-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 0.50 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether<	106-93-4	1,2-Dibromoethane (EDB)	ND	0.50
541-73-1 1,3-Dichlorobenzene ND 1.0 106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	74-95-3	Dibromomethane	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	95-50-1	1,2-Dichlorobenzene	ND	1.0
106-46-7 1,4-Dichlorobenzene ND 1.0 75-71-8 Dichlorodifluoromethane ND 1.0 75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	541-73-1	1,3-Dichlorobenzene	ND	1.0
75-34-3 1,1-Dichloroethane ND 1.0 107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	106-46-7	1,4-Dichlorobenzene	ND	1.0
107-06-2 1,2-Dichloroethane ND 1.0 75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-71-8	Dichlorodifluoromethane	ND	1.0
75-35-4 1,1-Dichloroethene ND 1.0 156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-34-3	1,1-Dichloroethane	ND	1.0
156-59-2 cis-1,2-Dichloroethene ND 1.0 156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	107-06-2	1,2-Dichloroethane	ND	1.0
156-60-5 trans-1,2-Dichloroethene ND 1.0 78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	75-35-4	1,1-Dichloroethene	ND	1.0
78-87-5 1,2-Dichloropropane ND 1.0 142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-59-2	cis-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	156-60-5	trans-1,2-Dichloroethene	ND	1.0
142-28-9 1,3-Dichloropropane ND 1.0 590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	78-87-5	1,2-Dichloropropane	ND	1.0
590-20-7 2,2-Dichloropropane ND 1.0 563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	142-28-9		ND	1.0
563-58-6 1,1-Dichloropropene ND 1.0 10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	590-20-7		ND	1.0
10061-01-5 cis-1,3-Dichloropropene ND 0.50 10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	563-58-6		ND	1.0
10061-02-6 trans-1,3-Dichloropropene ND 0.50 60-29-7 Diethyl ether ND 1.0	10061-01-5		ND	0.50
60-29-7 Diethyl ether ND 1.0	10061-02-6		ND	0.50
·	60-29-7		ND	1.0
	123-91-1	•	ND	20

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Matrix: Aqueous

Percent Moisture: N/A

Customer: Fuss & O'Neill Laboratory: Premier Laboratory, LLC Location: Franklin, MA

Project: 20050458.B10/Nu-Style Phase II PL Report No: E711585 PL Sample No: 12 (continued) Sample Description: 841071107-12

Date Collected: 11/6/2007 Date Received: 11/8/2007 Date Extracted: By: Date Analyzed: 11/13/07 By: DDD

Sample Weight/Volume: Dilution Factor: 1 Method: 8260B Soil Extract Volume: QC Batch#: 57819 Lab Data File: J32805.D

Units: ug/L

CAS No.	Parameter	Re	esult	DL
100-41-4	Ethylbenzene		ND	1.0
87-68-3	Hexachlorobutadiene		ND	0.50
591-78-6	2-Hexanone		ND	5.0
98-82-8	Isopropylbenzene		ND	1.0
99-87-6	4-Isopropyltoluene		ND	1.0
1634-04-4	Methyl tert-butyl ether (MTBE)		ND	1.0
108-10-1	4-Methyl-2-pentanone (MIBK)		ND	5.0
75-09-2	Methylene chloride		ND	5.0
91-20-3	Naphthalene		ND	1.0
103-65-1	n-Propylbenzene		ND	1.0
100-42-5	Styrene		ND	1.0
109-99-9	Tetrahydrofuran		ND	1.0
110-57-6	trans-1,4-Dichloro-2-butene		ND	5.0
76-13-1	1,1,2-Trichloro-1,2,2-trifluoroethane		ND	1.0
96-18-4	1,2,3-Trichloropropane		ND	1.0
630-20-6	1,1,1,2-Tetrachloroethane		ND	1.0
79-34-5	1,1,2,2-Tetrachloroethane		ND	0.50
127-18-4	Tetrachloroethene (PCE)		ND	1.0
108-88-3	Toluene		ND	1.0
87-61-6	1,2,3-Trichlorobenzene		ND	1.0
120-82-1	1,2,4-Trichlorobenzene		ND	1.0
71-55-6	1,1,1-Trichloroethane		ND	1.0
79-00-5	1,1,2-Trichloroethane		ND	1.0
79-01-6	Trichloroethene (TCE)		ND	1.0
75-69-4	Trichlorofluoromethane		ND	1.0
95-63-6	1,2,4-Trimethylbenzene		ND	1.0
108-67-8	1,3,5-Trimethylbenzene		ND	1.0
75-01-4	Vinyl chloride		ND	1.0
95-47-6	o-Xylene		ND	1.0
108-38-3	m,p-Xylenes		ND	1.0
Surrogate	Recovery	Limits		
1,2-Dichloroethane-d4	103%	89%-113%		
_ ~ .				

Surrogate	Recovery	Limits
1,2-Dichloroethane-d4	103%	89%-113%
Bromofluorobenzene	89%	83%-107%
Toluene-d8	98%	88%-108%

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	1	Sample Description:	841071106-01
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Extracted:	11/12/07	Percent Moisture:	N/A
Date Analyzed:	-11/12/07	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	78	40%-140%
2-Bromonaphthalene	NA	40%-140%
2-Fluorobiphenyl	NA	40%-140%
o-Terphenyl	116	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	2	Sample Description:	841071106-02
Preservative	HCL	•	
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Extracted:	11/12/07	Percent Moisture:	N/A
Date Analyzed:	-11/12/07	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	66	40%-140%
2-Bromonaphthalene	NA	40%-140%
2-Fluorobiphenyl	NA	40%-140%
o-Terphenyl	94	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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EXTRACTABLE PETROLEUM HYDROCARBON (EPH)

Laboratory:	Premier Laboratory, LLC	Client: Location:	Fuss & O'Neill Franklin, MA
PL Report No:	E711585	Project:	20050458.B10/Nu-Style Phase II
PL Sample No:	3	Sample Description:	841071106-03
Preservative	HCL		
		Dilution (Target):	1
Date Collected:	11/6/2007		
Date Received:	11/8/2007	Matrix:	Aqueous
Date Extracted:	11/12/07	Percent Moisture:	N/A
Date Analyzed:	-11/12/07	Method:	MADEP EPH
		Ext Method:	3545

(EPH) RANGE RESULTS

Parameter	Parameter Dilution	Results	QL	Units
C9-C18 Aliphatics	1	ND	100	ug/L
C19-C36 Aliphatics	1	ND	100	ug/L
C11-C22 Aromatics*	1	ND	100	ug/L

^{*} Excludes Targeted PAH Analytes

SURROGATE RECOVERIES

Surrogate	%Recovery	Acceptance Range
1-Chlorooctadecane	69	40%-140%
2-Bromonaphthalene	NA	40%-140%
2-Fluorobiphenyl	NA	40%-140%
o-Terphenyl	104	40%-140%

TARGETED PAH ANALYTES

Analyte	Results	QL	Units
2-Methylnaphthalene	ND	1.0	ug/L
Acenaphthene	ND	1.0	ug/L
Acenaphthylene	ND	1.0	ug/L
Anthracene	ND	1.0	ug/L
Benzo[a]anthracene	ND	1.0	ug/L
Benzo[a]pyrene	ND	0.20	ug/L
Benzo[b]fluoranthene	ND	1.0	ug/L
Benzo[g,h,i]perylene	ND	0.50	ug/L
Benzo[k]fluoranthene	ND	1.0	ug/L
Chrysene	ND	1.0	ug/L
Dibenz[a,h]anthracene	ND	0.50	ug/L
Fluoranthene	ND	1.0	ug/L
Fluorene	ND	1.0	ug/L
Indeno[1,2,3-cd]pyrene	ND	0.50	ug/L
Naphthalene	ND	1.0	ug/L
Phenanthrene	ND	1.0	ug/L
Pyrene	ND	1.0	ug/L

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Lab Name: Premier Laboratory, LLC Date Analyzed: 11/13/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1113 Location: Franklin, MA

Lab File ID: J32799.D

	Spike	Sample		QC
	Added.	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec #	Rec
1,1,1,2-Tetrachloroethane	50.00	51.91	104	79-115
1,1,1-Trichloroethane	50.00	56.37	113	77-123
1,1,2,2-Tetrachloroethane	50.00	48.58	97	72-120
1,1,2-Trichloroethane	50.00	52.59	105	80-114
1,1-Dichloroethane	50.00	52.08	104	80-114
1,1-Dichloroethene	50.00	53.40	107	75-128
1,1-Dichloropropene	50.00	54.77	110	75-117
1,2,3-Trichlorobenzene	50.00	57.60	115	70-127
1,2,4-Trichlorobenzene	50.00	56.08	112	70-123
1,2,4-Trimethylbenzene	50.00	49.42	99	76-126
1,2-Dibromoethane (EDB)	50.00	53.76	108	82-118
1,2-Dichlorobenzene	50.00	48.74	97	76-121
1,2-Dichloroethane	50.00	52.78	106	75-115
1,2-Dichloropropane	50.00	53.83	108	71-130
1,3,5-Trimethylbenzene	50.00	52.54	105	76-122
1,3-Dichlorobenzene	50.00	47.88	96	79-120
1,3-Dichloropropane	50.00	53.66	107	82-118
1,4-Dichlorobenzene	50.00	47.00	94	77-116
1,4-Dioxane	50.00	52.06	104	70-119
2,2-Dichloropropane	50.00	58.48	117	71-125
2-Butanone (MEK)	50.00	43.24	86	70-128
4-Chlorotoluene	50.00	48.53	97	77-117
4-Isopropyltoluene	50.00	47.06	94	78-124
4-Methyl-2-pentanone (MIBK)	50.00	40.64	81	75-121
Acetone	50.00	46.56	93	70-118
Benzene	50.00	53.41	107	77-118
Bromobenzene	50.00	49.86	100	79-116
Bromochloromethane	50.00	56.85	114	80-122

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/13/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1113 Location: Franklin, MA

Lab File ID: J32799.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Campound	(ug/L)	(ug/L)	Rec #	Rec
Bromodichloromethane	50.00	52.84	106	82-127
Bromoform	50.00	47.79	96	78-122
Bromomethane	50.00	49.13	98	70-130
Carbon disulfide	50.00	44.61	89	70-130
Carbon tetrachloride	50.00	49.86	100	77-125
Chlorobenzene	50.00	51.05	102	80-118
Chloroform	50.00	53.55	107	80-113
Chloromethane	50.00	44.36	89	70-130
cis-1,2-Dichloroethene	50.00	53.68	107	85-120
cis-1,3-Dichloropropene	50.00	53.26	106	79-116
Di-isopropyl ether (DIPE)	50.00	44.12	88	78-121
Dibromochloromethane	50.00	51.74	103	79-122
Dibromomethane	50.00	55.36	111	78-120
Ethyl tertiary-butyl ether (EtBE)	50.00	40.66	81	81-122
Ethylbenzene	50.00	55.16	110	84-123
Hexachlorobutadiene	50.00	46.71	93	70-121
Isopropylbenzene	50.00	52.30	105	78-120
m,p-Xylenes	100.0	108.8	109	75-129
Methyl tert-butyl ether (MTBE)	50.00	46.32	93	70-127
Methylene chloride	50.00	48.66	97	72-128
n-Butylbenzene	50.00	51.38	103	70-124
n-Propylbenzene	50.00	53.08	106	80-127
Naphthalene	50.00	54.02	108	70-126
o-Xylene	50.00	51.22	102	78-118
sec-Butylbenzene	50.00	48.99	98	78-118
Styrene	50.00	50.60	101	83-124
tert-Butylbenzene	50.00	51.64	103	76-118
Tertiary-amyl methyl ether (TAME)	50.00	42.46	85	81-123

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/13/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1113 Location: Franklin, MA

Lab File ID: J32799.D (continued)

	Spike	Sample		QC
	Added.	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec #	Rec
Tetrachloroethene (PCE)	50.00	52.86	106	77-122
Toluene	50.00	53.03	106	78-120
trans-1,2-Dichloroethene	50.00	51.41	103	80-120
trans-1,3-Dichloropropene	50.00	45.16	90	71-111
Trichloroethene (TCE)	50.00	54.97	110	74-118
Vinyl chloride	50.00	46.46	93	70-116

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/14/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1114 Location: Franklin, MA

Lab File ID: J32823.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Campound	(ug/L)	(ug/L)	Rec #	Rec
1,1,1,2-Tetrachloroethane	50.00	52.23	104	79-115
1,1,1-Trichloroethane	50.00	52.41	105	77-123
1,1,2,2-Tetrachloroethane	50.00	51.17	102	72-120
1,1,2-Trichloroethane	50.00	52.61	105	80-114
1,1-Dichloroethane	50.00	49.18	98	80-114
1,1-Dichloroethene	50.00	53.70	107	75-128
1,1-Dichloropropene	50.00	50.68	101	75-117
1,2,3-Trichlorobenzene	50.00	59.88	120	70-127
1,2,4-Trichlorobenzene	50.00	57.79	116	70-123
1,2,4-Trimethylbenzene	50.00	49.10	98	76-126
1,2-Dibromoethane (EDB)	50.00	51.65	103	82-118
1,2-Dichlorobenzene	50.00	51.06	102	76-121
1,2-Dichloroethane	50.00	50.83	102	75-115
1,2-Dichloropropane	50.00	50.65	101	71-130
1,3,5-Trimethylbenzene	50.00	50.44	101	76-122
1,3-Dichlorobenzene	50.00	49.28	98	79-120
1,3-Dichloropropane	50.00	52.48	105	82-118
1,4-Dichlorobenzene	50.00	48.90	98	77-116
1,4-Dioxane	50.00	52.73	105	70-119
2,2-Dichloropropane	50.00	56.22	112	71-125
2-Butanone (MEK)	50.00	49.02	98	70-128
4-Chlorotoluene	50.00	48.34	97	77-117
4-Isopropyltoluene	50.00	49.52	99	78-124
4-Methyl-2-pentanone (MIBK)	50.00	45.16	90	75-121
Acetone	50.00	47.94	96	70-118
Benzene	50.00	48.76	98	77-118
Bromobenzene	50.00	49.47	99	79-116
Bromochloromethane	50.00	52.95	106	80-122

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/14/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1114 Location: Franklin, MA

Lab File ID: J32823.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Campound	(ug/L)	(ug/L)	Rec #	Rec
Bromodichloromethane	50.00	51.82	104	82-127
Bromoform	50.00	50.98	102	78-122
Bromomethane	50.00	51.64	103	70-130
Carbon disulfide	50.00	47.41	95	70-130
Carbon tetrachloride	50.00	55.70	111	77-125
Chlorobenzene	50.00	50.04	100	80-118
Chloroform	50.00	50.00	100	80-113
Chloromethane	50.00	47.30	95	70-130
cis-1,2-Dichloroethene	50.00	51.44	103	85-120
cis-1,3-Dichloropropene	50.00	51.10	102	79-116
Di-isopropyl ether (DIPE)	50.00	49.54	99	78-121
Dibromochloromethane	50.00	53.17	106	79-122
Dibromomethane	50.00	52.08	104	78-120
Ethyl tertiary-butyl ether (EtBE)	50.00	45.85	92	81-122
Ethylbenzene	50.00	52.29	104	84-123
Hexachlorobutadiene	50.00	47.69	95	70-121
Isopropylbenzene	50.00	50.14	100	78-120
m,p-Xylenes	100.0	105.6	106	75-129
Methyl tert-butyl ether (MTBE)	50.00	50.72	101	70-127
Methylene chloride	50.00	46.78	94	72-128
n-Butylbenzene	50.00	51.81	104	70-124
n-Propylbenzene	50.00	51.38	103	80-127
Naphthalene	50.00	57.64	115	70-126
o-Xylene	50.00	51.56	103	78-118
sec-Butylbenzene	50.00	50.96	102	78-118
Styrene	50.00	50.50	101	83-124
tert-Butylbenzene	50.00	51.80	104	76-118
Tertiary-amyl methyl ether (TAME)	50.00	44.14	88	81-123

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/14/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1114 Location: Franklin, MA

Lab File ID: J32823.D (continued)

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec #	Rec
Tetrachloroethene (PCE)	50.00	50.30	101	77-122
Toluene	50.00	49.44	99	78-120
trans-1,2-Dichloroethene	50.00	51.53	103	80-120
trans-1,3-Dichloropropene	50.00	47.24	94	71-111
Trichloroethene (TCE)	50.00	51.68	103	74-118
Vinyl chloride	50.00	49.81	100	70-116

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 4 8260B Method Blank Summary

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Lab File ID: J32802.D Lab Sample ID: VBLK1113

Matrix: Water Date Analyzed: 11/13/07

Instrument ID: MS8 Date Extracted:

Time Analyzed: 1112

This Method Blank Applies To The Following Samples, MD and MSD:

	Iab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711585-1A	841071106-01	J32815.D	11/13/2007
2	E711585-2A	841071106-02	J32816.D	11/13/2007
3	E711585-3A	841071106-03	J32817.D	11/13/2007
4	E711585-4A	841071106-04	J32818.D	11/13/2007
5	E711585-5A	841071106-05	J32819.D	11/13/2007
6	E711585-7	841071106-07	J32804.D	11/13/2007
7	E711585-12	841071107-12	J32805.D	11/13/2007
8	VLCS1113	VLCS1113	J32799.D	11/13/2007
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FORM 4 8260B Method Blank Summary

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Lab File ID: J32826.D Lab Sample ID: VBLK1114

Matrix: Water Date Analyzed: 11/14/07

Instrument ID: MS8 Date Extracted:

Time Analyzed: 1103

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711585-6A	841071106-06	J32827.D	11/14/2007
2	E711585-8A	841071107-08	J32828.D	11/14/2007
3	E711585-9A	841071107-09	J32829.D	11/14/2007
4	E711585-10A	841071107-10	J32830.D	11/14/2007
5	E711585-11A	841071107-11	J32831.D	11/14/2007
6	VLCS1114	VLCS1114	J32823.D	11/14/2007
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FORM 3 Water MADEP VPH Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/16/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1116 Location: Franklin, MA

Lab File ID: 2111623.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec #	Rec
Benzene	25.00	24.67	99	70-130
Ethylbenzene	25.00	24.51	98	70-130
m,p-Xylenes	50.00	48.09	96	70-130
Methyl tert-butyl ether (MTBE)	25.00	23.49	94	70-130
Naphthalene	25.00	21.57	86	70-130
o-Xylene	25.00	23.82	95	70-130
Toluene	25.00	24.10	96	70-130
Benzene #2	25.00	23.55	94	70-130
Ethylbenzene #2	25.00	24.21	97	70-130
Naphthalene #2	25.00	21.58	86	70-130
Toluene #2	25.00	23.77	95	70-130

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Water MADEP VPH Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/16/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: VLCS1116 DUP Location: Franklin, MA

Lab File ID: 2111634.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Compound	(ug/L)	(ug/L)	Rec #	Rec
Benzene	25.00	23.71	95	70-130
Ethylbenzene	25.00	23.89	96	70-130
m,p-Xylenes	50.00	46.67	93	70-130
Methyl tert-butyl ether (MTBE)	25.00	23.34	93	70-130
Naphthalene	25.00	23.19	93	70-130
o-Xylene	25.00	23.14	92	70-130
Toluene	25.00	23.32	93	70-130
Benzene #2	25.00	24.23	97	70-130
Ethylbenzene #2	25.00	23.85	95	70-130
Naphthalene #2	25.00	23.10	92	70-130
Toluene #2	25.00	23.71	95	70-130

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 4 MADEP VPH Method Blank Summary

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Lab File ID: 2111622.D Lab Sample ID: VBLK1116

Matrix: Water Date Analyzed: 11/16/07

Instrument ID: GC2 Date Extracted:

Time Analyzed: 1631

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711585-1A	841071106-01	2111625.D	11/16/2007
2	E711585-2A	841071106-02	2111626.D	11/16/2007
3	E711585-3A	841071106-03	2111627.D	11/16/2007
4	VLCS1116	VLCS1116	2111623.D	11/16/2007
5	VLCS1116 DUP	VLCS1116 DUP	2111634.D	11/16/2007
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FORM 3 Water MADEP EPH Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/12/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1112A-1 Location: Franklin, MA

Lab File ID: A23912.D

	Spike	Sample		QC
	Added	Concentration	%	Limits
Campound	()	()	Rec #	Rec
2-Methylnaphthalene	40.00	34.30	86	40-140
Acenaphthene	40.00	32.28	81	40-140
Acenaphthylene	40.00	32.67	82	40-140
Anthracene	40.00	38.44	96	40-140
Benzo[a]anthracene	40.00	35.43	88	40-140
Benzo[a]pyrene	40.00	35.89	90	40-140
Benzo[b]fluoranthene	40.00	35.20	88	40-140
Benzo[g,h,i]perylene	40.00	41.55	104	40-140
Benzo[k]fluoranthene	40.00	34.60	86	40-140
Chrysene	40.00	33.86	85	40-140
Decane	40.00	38.21	96	40-140
Dibenz[a,h]anthracene	40.00	39.53	99	40-140
Docosane	40.00	42.90	107	40-140
Dodecane	40.00	40.94	102	40-140
Eicosane	40.00	43.01	108	40-140
Fluoranthene	40.00	34.25	86	40-140
Fluorene	40.00	32.83	82	40-140
Hexacosane	40.00	41.85	105	40-140
Hexadecane	40.00	40.15	100	40-140
Hexatriacontane	40.00	56	140	40-140
Indeno[1,2,3-cd]pyrene	40.00	40.08	100	40-140
Naphthalene	40.00	34.52	86	40-140
Nonadecane	40.00	42.87	107	40-140
Nonane	40.00	37.03	92	30-140
Octacosane	40.00	40.95	102	40-140
Octadecane	40.00	43.85	110	40-140
Phenanthrene	40.00	29.67	74	40-140
Pyrene	40.00	33.84	85	40-140

[#] Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Water MADEP EPH Lab Control Sample

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/12/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1112A-1 Location: Franklin, MA

Lab File ID: A23912.D (continued)

	Spike	Sample		QC
	Added.	Concentration	%	Limits
Campound	()	()	Rec #	Rec
Tetracosane	40.00	42.25	106	40-140
Tetradecane	40.00	39.93	100	40-140
Triacontane	40.00	44.34	111	40-140

Column to be used to flag recovery values with an asterisk

^{*} Values outside of QC limits

FORM 3 Water MADEP EPH Lab Control Sample Duplicate

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/12/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: ICS1112A-1 Location: Franklin, MA

Lab File ID: A23913.D

	Spike	Sample				QC
	Added	Concentration	%		L	imits
Campound	()	()	Rec #	RPD#	RPD	Rec
2-Methylnaphthalene	40.00	28.86	72	17.7	25	40-140
Acenaphthene	40.00	27.77	69	16.0	25	40-140
Acenaphthylene	40.00	27.62	69	17.2	25	40-140
Anthracene	40.00	33.10	83	14.5	25	40-140
Benzo[a]anthracene	40.00	32.01	80	9.52	25	40-140
Benzo[a]pyrene	40.00	33.18	83	8.09	25	40-140
Benzo[b]fluoranthene	40.00	31.86	80	9.52	25	40-140
Benzo[g,h,i]perylene	40.00	37.88	95	9.04	25	40-140
Benzo[k]fluoranthene	40.00	31.30	78	9.76	25	40-140
Chrysene	40.00	30.53	76	11.2	25	40-140
Decane	40.00	32.28	81	16.9	25	40-140
Dibenz[a,h]anthracene	40.00	35.97	90	9.52	25	40-140
Docosane	40.00	38.46	96	10.8	25	40-140
Dodecane	40.00	34.19	85	18.2	25	40-140
Eicosane	40.00	37.83	94	13.9	25	40-140
Fluoranthene	40.00	30.15	75	13.7	25	40-140
Fluorene	40.00	27.95	70	15.8	25	40-140
Hexacosane	40.00	38.16	95	10.0	25	40-140
Hexadecane	40.00	34.21	86	15.0	25	40-140
Hexatriacontane	40.00	51.45	129	10.3	25	40-140
Indeno[1,2,3-cd]pyrene	40.00	36.34	91	9.42	25	40-140
Naphthalene	40.00	29.02	72	17.7	25	40-140
Nonadecane	40.00	37.46	94	12.9	25	40-140
Nonane	40.00	33.02	82	11.5	25	30-140
Octacosane	40.00	38.31	96	6.06	25	40-140
Octadecane	40.00	37.62	94	15.7	25	40-140
Phenanthrene	40.00	30.10	75	1.34	25	40-140
Pyrene	40.00	29.98	75	12.5	25	40-140

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 3 Water MADEP EPH Lab Control Sample Duplicate

Lab Name: Premier Laboratory, LLC Date Analyzed: 11/12/07

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Sample No.: LCS1112A-1 Location: Franklin, MA

Lab File ID: A23913.D (continued)

	Spike	Sample				QC
	Added	Concentration	%		L	imits
Compound	()	()	Rec #	RPD#	RPD	Rec
Tetracosane	40.00	38.39	96	9.90	25	40-140
Tetradecane	40.00	33.63	84	17.4	25	40-140
Triacontane	40.00	40.97	102	8.45	25	40-140

[#] Column to be used to flag recovery and RPD values with an asterisk

^{*} Values outside of QC limits

FORM 4 MADEP EPH Method Blank Summary

Project No.: E711585 Project: 20050458.B10/Nu-Style Phase II

Lab File ID: A23914.D Lab Sample ID: E1112BA-1

Matrix: Water Date Analyzed: 11/12/07

Instrument ID: GCl Date Extracted:

Time Analyzed: 1124

This Method Blank Applies To The Following Samples, MD and MSD:

	Lab	Client	Lab	Date
	Sample No.	Sample ID	File ID	Analyzed
1	E711585-1C	841071106-01	A23917.D	11/12/2007
2	E711585-2C	841071106-02	A23918.D	11/12/2007
3	E711585-3C	841071106-03	A23919.D	11/12/2007
4	LCS1112A-1	ICS1112A-1	A23912.D	11/12/2007
5	LCSD1112A-1	LCSD1112A-1	A23913.D	11/12/2007
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146 Hartford Road, Manchester, CT 06040 56 Quarry Road, Trumbull, CT 06611

1419 Richland Street, Columbia, SC 29201

78 Interstate Drive, West Springfield, MA 01089

1 610 Lynndale Court, Suite E, Greenville, NC 27858

☐ 24 Madison Avenue Extension, Albany, NY 12203

275 Promenade Street, Suite 350, Providence, RI 02908

30 Washington Street, Suite 301, Poughkeepsie, NY 12601

Other_

CHAIN-OF-CUST	DDY RECORD	14728	☐ 1 Day* ☐ 3 Days* ☐ 2 Days* ☐ 2 Days* ☐ Standard (days)	Char (days)
Project Name	PROJECT LOCATION	Project Number	- Lucio	LABORATORY
Nu-Style Physe II	Franklin, MA	20050458.B	10	Premier
REPORT TO: Double foss		Analysis ///		Containers
INVOICE TO: DOLAND FOSS		Request		\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\\
P.O. No.: 84120050458BID		- / \J : 5 / \J		
Sampler's Signature:	Date: // 1/0-			/ /:>/:>/:>/ //
Source Codes: MW=Monitoring Well SW=Surface Water T=Treatment Facility S=Soil B=Sediment X=Other T s Slank	W=Waste A=Air		/	
Item No. 1 2 3 4 Sample Number	Source Date Time Code Sampled Sample			Comments
1 / 84107-1106-01	mw 11/6/07 100		421	
2 1 -02	100		421	
3 -03	_ 100	V V V	421	
4 -04	134		2 1	
5 -05	142		2 1	
c -06	153		2 1	
7 / -07	V 160			Trig Blank
8 84107-1107-08	MW 11/7/07 WZ		S2	
9 109	1 1 1100		2	1 Feld-Sitead
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Transfer Relinquished By	Accepted By D	SEE ATTAMA	quirements: <u>ED_BBn 7.0194E_Fibre_R</u>	7/4
1 / h Bu For	Shalge "16	Additional Comments:		
3 Tim May Jungle	J_ 5	16 0495 PLEASE Con	PLEASE ATTATCHED C	HECKLIST
4 1/2 2 1/4/11/1/	SMACA.	N SE		



☐ 146 Hartford Road, Manchester, CT 06040 1 56 Quarry Road, Trumbell, CT 06611

☐ 1419 Richland Street, Columbia, SC 29201

3 78 Interstate Drive, West Springfield, MA	01089
☐ 610 Lynndale Court, Suite E, Greenville, N	C 27858

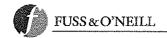
☐ 24 Madison Avenue Extension, Albany, NY 12203

275 Promenade Street, Suite 350, Providence, RI-02908

3 80 Washington Street, Suite 301, Poughkeepsie, NY 12601

Other___

CHAIN-OF-CUST	TODY RECORD	14729	Turnaround 1 1 Day* 1 3 Days* 2 2 Days* X Standard (1 Other (days)
PROJECT NAME Nu-Style Phase II	PROJECT LOCATION Frankling	Project Number Quesoyss , C		LABORATORY Premyer
REPORT TO: DONAL FOST		Analysis / /		Containers
INVOICE TO: David Foss		Request		
P.O. No.: 841 20050458 B10				
Sampler's Signature:	Date: 1/1/2/17			/ \/ \/ \&/ \/ /
Source Codes: MW=Monitoring Well PW=Potable Water S=Soil SW=Surface Water T=Treatment Facility B=Sediment K=Other Blank	W=Waste			
Item No. 1 2 3 4 Sample Number	Source Date Time Code Sampled Sampled			Comments
11 84107407-11	MW 11/7/07 1145		2 1	Feld Fiteel
2 -12	<u> </u>	√		TopBlace
Transfer Number Relinquished By	Accepted By Da	SEE ATTAR	uirements: MED TABLE FOR RL	
1 Man Bour	o fielge 1/5	1330 Additional Comments:	The state of the s	
2 1 5 - Sily 12 12	the way	100 0945 Please Con	ylete attacked	
4 Parlliant	11-6			cuck/13t



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

			<u>YES</u>	NO COMMENTS
1.	SDG Project Narratives	K		**************************************
2.	Traffic Report			<u> VIA</u>
3.	Volatiles Data			
	a. Sample Data			
	Target Compound List (TCL) Results			
	Reconstructed total ion chromatograms (RIC) for each			
	Sample	Ŋ.		
	For each sample:	c		
	Raw spectra and background-subtracted mass spectra			
	target compounds identified	<u>⊠</u> .	Ц	
	Mass spectra of all reported TICs with three best library	•		NIA
	Percent solids calculations			VIA
	1 CICCIE SONGS CAICHANOIIS		لــا	
	b. Standards Data (all instruments)			
	Initial Calibration Data	\bar{\bar{\bar{\bar{\bar{\bar{\bar{		
	RICs and Quan Reports for all Standards	∑ -		
	Continuing Calibration			
	RICs and Quan Reports for all Standards	M		
	Internal Standard Area Summary	A		
	c. Raw QC Data			
	Blank Data			
	Matrix Spike Data			MIA
	Matrix Spike Duplicate Data			w/A
4.	Semivolatiles Data			
	a. QC Summary			
	Surrogate Percent Recovery Summary			
	MS/MSD Summary			NIA
	Method Blank Summary	<u>M</u> -		
	Tuning and Mass Calibration			NIA



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTSP**

		<u>YES</u>	<u>NO</u>	<u>COMMENTS</u>
b.	Sample Data TCL Results	[]		
	Tentatively Identified Compounds			MA
	Reconstructed total ion chromatograms (RIC) for each Sample			NIA
	For each sample: Raw spectra and background-subtracted mass			
	spectra of TCL compounds			NIA
	Mass spectra of TICs with 3 best library matches			NIA
	GPC chromatograms (if GPC performed)			NIA
c.	Standards Data (all instruments)			
	Initial Calibration Data			
	RICs and Quan Reports for all Standards Continuing Calibration	<u>₩</u>		
	RICs and Quan Reports for all Standards	K N		
	Internal Standard Areas Summary			MA
	Internal Standard Areas Summary			NIA
d.	Raw QC Data			
	Decafluorotripbenylphosphine (DFTPP)			MIA
	Blank Data			
	Matrix Spike Data			NIA
	Matrix Spike Duplicate Data		ᆸ _	WIA
M	scellaneous Data			
	Original preparation and analysis forms or copies of preparation			
	and analysis log book pages	[2]-	□ .	
	Internal sample & sample extract transfer chain-of custody records			
	Screening Records			WIA
	All instrument output, including strip charts from screening			(** ())
	activities (describe or list)	_ _	_	



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST ORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

		$\underline{\text{YES}}$	<u>NO</u>	COMMENTS	<u> </u>
6.	Chain-of-Custody Records Sample Log-in Sheet (Lab & DC1) Miscellaneous Shipping/Receiving Records (describe or	⊠ □ list) □		NIA NIA	
7.	Internal Lab Sample Transfer Records and Tracking Sheets (describe or list)		□		
8.	Other Records (describe or list)				
9.	Comments:				
	** See laboratory Quality Assurance Plan for limits.				
	Completed by: Signature) Completed by: Signature Completed by: Signat	l Onten C Mana Title)	ics ger		9/07
	I certify that the above information is true and accurate. I further certified above analyses will be made available for review for seven (7) years.				
	Certified by: Lab (Signature) Robert Stee (Printed Name/1	renson/ba Pitle)	o Dire	ctor 11-19 Date	-07



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?***

	7	ES	NO	O COMMENTS
SDG Project Narratives	Ī	2		
Inorganic Analysis Data Sheet	A.	7		
Initial and Continuing Calibration Verification	Ē	Ź		
CRDL Standard for AA and ICP	Ĺ	đ		
Blanks		ZÍ		
ICP Interference Check Sample	ľ	7		
Spike Sample Recovery	/\bar{L}	9		
Post Digest Spike Sample Recovery	E			MA
Duplicates	_/_]		-
Laboratory Control Sample	72			
Standard Addition Results]	Ø	NA
ICP Serial Dilutions	Z	Ī		
Instrument Detection Limits, Quarterly				
ICP Interelement Correction Factors, Annually	V	Í		
ICP Linear Ranges Quarterly				\
Preparation Log	$I_{\!$			
Analysis Run Log	V	ĺ		
ICP Raw Data	, V	j		
Furnace AA Raw Data	Е]	Į.	VA
Mercury Raw Data	D	Ì		
Percent Solids Calculations]	vÍ	
Digestion Logs	V			
EPA Shipping/Receiving Records				
(List all individual records)				
Chain-of Custody Records	Ţ			
Sample Log-In sheet	Č			
Miscellaneous Shipping/Receiving Records				
(List all individual records)				



PHASE II SITE ASSESSMENT FORMER NU-STYLE COMPANY, INC. FACILITY LABORATORY MODIFIED TIER II DATA VALIDATION CHECKLIST INORGANIC COMPOUNDS (Continued)

PERFORMED AND, WHERE APPLICABLE, WITHIN ACCEPTABLE LIMTS?**

YES NO COMMENTS

25.		al Lab Sample Transfer Re ribe or List) 1000004-	cords and Tracking Sheets			
26.	(Descri Prepar Analys Descri		•		□ □ <u>logbo</u> □ <u>hovdu</u>	ok opy + ULdronic-
27.	Other	Records (Describe or List))	_		
28.	Comm	nents:			······································	
**	See lab	poratory Quality Assurance	Plan for limits.		·	
Compl (Lab)	leted by:	B. Truppy wyn fri (Signature)	BOY <u>boyo Supryuyuski</u> (Printed Name/Title	Jno.	ganie Hov	11/19/04 Date
associa	ited with		e and accurate. I further cert made available for review fo	•		-
Certific (Lab)	ed by:	Manager (Signature)	Robert Stevenson (Printed Name/Title	1646	Ducker	11-14-07 Date



APPENDIX E

Copy of BWSC107 Transmittal Forms



TIER CLASSIFICATION TRANSMITTAL FORM Pursuant to 310 CMR 40.0500 (Subpart E)

R	Λ	19	C1	07	
D	٧v	0	u	U1	

Release Tracking Number

	1 disdant to 310 Civil (40.0300 (Subpart L)
A. D	ISPOSAL SITE LOCATION:
1. D	isposal Site Name:
2 S	treet Address:
2. 0	ilicot / iddi coss.
3. C	ity/Town: 4. ZIP Code:
B. T	HIS FORM IS BEING USED TO: (check all that apply)
	1. Submit a new Tier Classification Submittal for a Tier I Site , including a Numerical Ranking Scoresheet (BWSC107A) (check one). A Tier I Permit Application must also be submitted.
	a. Tier IA b. Tier IB c. Tier IC
	2. Submit a new Tier Classification Submittal for a Tier II Site , including the Numerical Ranking Scoresheet (BWSC107A) and the Tier II Compliance History (BWSC107B).
	3. Submit a Phase I Completion Statement as per 310 CMR 40.0480.
	If previously submitted, provide date mm/dd/yyyy
	4. Submit a Phase II Scope of Work as per 310 CMR 40.0834.
	If previously submitted, provide datemmm/dd/yyyy
	5. Submit a Phase II Conceptual Scope of Work supporting a Tier Classification Submittal.
	6. Submit a Tier II Extension Submittal for Response Actions at a Tier II Site including the Tier II Compliance History (BWSC107B).
	7. Submit a Tier II Transfer Submittal for a change in person(s) undertaking Response Actions at a Tier II Site including the Tier II Compliance History (BWSC107B) and the Tier II Transferor Certification (BWSC107C).
	Proposed effective date of transfer : mm/dd/yyyy
	8. Submit a Revised Tier Classification Submittal, including a Numerical Ranking Scoresheet (BWSC107A). A Major Permit Modification may also need to be submitted.
	If this revised submittal is re-classifying the site check the new classification.
	a. Tier IA b. Tier IB c. Tier IC d. Tier II
	9. Submit a Notice that an additional Release Tracking Number(s) is (are) being linked to this Tier Classified Site (Primary RTN). Future response actions addressing the Release or Threat of Release notification condition associated with additional Release Tracking Numbers (RTNs) will be conducted as part of the Response Actions planned or ongoing at the Primary Site listed above. For a previously Tier Classified Primary Site, if there is a reasonable likelihood that the addition of the new secondary RTN(s) would change the classification of the site, a Revised Tier Classification Submittal must also be made.
	Provide Release Tracking Number(s): a b
	All future Response Actions must occur according to the deadlines applicable to the Primary RTN. Use only the Primary RTN when making future submittals for this site unless specifically relating to response actions started before the linking occurred.

Revised: 02/10/2004 Page 1 of 4



Massachusetts Department of Environmental Protection

Bureau of Waste Site Cleanup

TIER CLASSIFICATION TRANSMITTAL FORM

Pursuant to 310 CMR 40.0500 (Subpart E)

D1100101							
Rele	ase T	racking Numb	er				
	_						

RWSC107

C. LSP SIGNATURE AND STAMP:

I attest under the pains and penalties of perjury that I have personally examined and am familiar with this transmittal form, including any and all documents accompanying this submittal. In my professional opinion and judgment based upon application of (i) the standard of care in 309 CMR 4.02(1), (ii) the applicable provisions of 309 CMR 4.02(2) and (3), and 309 CMR4.03(2), and (iii) the provisions of 309 CMR 4.03(3), to the best of my knowledge, information and belief,

- > if Section B of this form indicates that a **Tier I or Tier II Classification Submittal** including the **Numerical Ranking System Scoresheet** is being submitted, this Tier Classification Submittal has been developed in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000 and, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed and implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) complies(y) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B of this form indicates that a **Phase I Completion Statement** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed and implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B of this form indicates that a **Phase II Scope of Work** is being submitted, the response action(s) that is (are) the subject of this submittal (i) has (have) been developed in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) comply(ies) with the identified provisions of all orders, permits, and approvals identified in this submittal;
- > if Section B of this form indicates that a **Tier II Extension Submittal** or a **Tier II Transfer Submittal** is being submitted, the response action(s) that is (are) the subject of this submittal (i) is (are) being implemented in accordance with the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, (ii) is (are) appropriate and reasonable to accomplish the purposes of such response action(s) as set forth in the applicable provisions of M.G.L. c. 21E and 310 CMR 40.0000, and (iii) complies(y) with the identified provisions of all orders, permits, and approvals identified in this submittal.

I am aware that significant penalties may result, including, but not limited to, possible fines and imprisonment, if I submit information which I know to be false, inaccurate or materially incomplete.

1.	LSP#:				
2.	First Name:		3.	Last Name:	
4.	Telephone:		5. Ext.:	6. FAX:	
7.	Signature: _	Submitted via e-DEP			
8.		mm/dd/yyyy		9. LSP Stamp:	

Revised: 02/10/2004 Page 2 of 4



TIER CLASSIFICATION TRANSMITTAL FORM

BW	SC	10	7
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Release Tracking Number

Pursuant to 310 CMR 40.0500 (Subpart E)
D. PERSON MAKING SUBMITTAL:
1. Check all that apply: a. change in contact name b. change of address c. change in the person undertaking response actions
2. Name of Organization:
3. Contact First Name: 4. Last Name:
5. Street: 6. Title:
7. City/Town: 8. State: 9. ZIP Code:
10. Telephone: 11. Ext.: 12. FAX:
E DELATIONELUD DE DEDEON MAZINE CUDMITTAL TO DICROCAL CITE.
E. RELATIONSHIP OF PERSON MAKING SUBMITTAL TO DISPOSAL SITE:
1. RP or PRP a. Owner b. Operator c. Generator d. Transporter
e. Other RP or PRP Specify:
2. Fiduciary, Secured Lender or Municipality with Exempt Status (as defined by M.G.L. c. 21E, s. 2)
3. Agency or Public Utility on a Right of Way (as defined by M.G.L. c. 21E, s. 5(j))
4. Any Other Person Making Submittal Specify Relationship:
F. REQUIRED ATTACHMENT AND SUBMITTALS:
1. Check here if the Response Action(s) on which this opinion is based, if any, are (were) subject to any order(s), permit(s) and/or approval(s) issued by DEP or EPA. If the box is checked, you MUST attach a statement identifying the applicable provisions thereof.
2. Check here to certify that the Chief Municipal Officer and the Local Board of Health have been notified of the submittal of any Phase Reports to DEP.
3. Check here to certify that a Legal Notice of a Tier Classification or Re-classification Submittal has been or will be made according to 310 CMR 40.1403, and a copy of the notice sent to DEP, the Chief Municipal Officer and the Local Board of Health.
4. For a Tier II Extension Submittal, check here to certify that a statement summarizing why a Permanent or Temporary Solution has not been achieved at the Disposal Site is attached.
5. For a Tier II Transfer Submittal, check here to certify that a statement summarizing the reasons for the proposed change in person(s) undertaking the Response Actions is attached. All Response Actions must be completed by the deadline applicable to the person who first filed either a Tier Classification Submittal for the Disposal Site or received a Waiver of Approvals.
6. Check here if any non-updatable information provided on this form is incorrect, e.g., Site Name or Street Address. Send corrections to the DEP Regional Office.
7. Check here to certify that the LSP Opinion containing the material facts, data, and other information is attached.

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BWSC1	07
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TIER CLASSIFICATION TRANSMITTAL FORM Pursuant to 310 CMR 40.0500 (Subpart E)	Release Tracking Number
G. CERTIFICATION OF PERSON MAKING SUBMITTAL:	
1. I,	ocuments accompanying this aining the information, the curate and complete, and (iii) submittal. I/the person or entity
If submitting a Tier II Classification, Extension or Transfer, I also attest under the pains and penaltic person(s) or entity(ies) on whose behalf this submittal is made has/have personally examined and requirements of M.G.L. c. 21E and 310 CMR 40.0000; (ii) based upon my inquiry of the/those Licen employed or engaged to render Professional Services for the disposal site which is the subject of t person(s) or entity(ies) on whose behalf this submittal is made, and my/that person's(s') or entity's estimated costs of necessary response actions, that/those person(s) or entity(ies) has/have the teability to proceed with response actions for such site in accordance with M.G.L. c. 21E, 310 CMR 40 requirements; and (iii) that I am fully authorized to make this attestation on behalf of the person(s) of this submittal. I/the person(s) or entity(ies) on whose behalf this submittal is made is aware of the 40.0172 for notifying the Department in the event that I/the person(s) or entity(ies) on whose behalf that it/they is/are unable to proceed with the necessary response actions.	d am/is familiar with the used Site Professional(s) this Transmittal Form and of the s(ies') understanding as to the chnical, financial and legal 0.0000 and other applicable or entity(ies) legally responsible the requirements in 310 CMR
2. By: Submitted via e-DEP 3. Title:	
Signature	
4. For: 5. Date:	
(Name of person or entity recorded in Section D)	mm/dd/yyyy
6. Check here if the address of the person providing certification is different from address reco	orded in Section D.
8. City/Town: 9. State: 10. Z	ZIP Code:
11. Telephone: 12. Ext.: 13. FAX:	
YOU ARE SUBJECT TO AN ANNUAL COMPLIANCE ASSURANCE FEE OF UP TO BILLABLE YEAR FOR THIS DISPOSAL SITE. YOU MUST LEGIBLY COMPLETE AL SECTIONS OF THIS FORM OR DEP MAY RETURN THE DOCUMENT AS INCOMPL SUBMIT AN INCOMPLETE FORM, YOU MAY BE PENALIZED FOR MISSING A REQUI	LL RELEVANT LETE. IF YOU
Date Stamp (DEP USE ONLY):	

Revised: 02/10/2004 Page 4 of 4



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Release	Tracking Number
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	ERICAL RAN nt to 310 CMR 40		, ,	SCORESH	EET Release	Tracking Number	
A. NRS SCORESHEET SUN	IMARY SECTION:						
1. Classification Submittal:	(check one)	a. Initial NR	RS Score	b. Revised	NRS Score		
2. Disposal Site Score:							
II III IV V VI Total							
3. Disposal Site Classificat	ion: (check one)						
a. Tier IA b	. Tier IB	c. Tier IC	d. Tier II				
B. DISPOSAL SITE INFORM	MATION (NRS SEC	CTION I):					
1. UTM Coordinates: a.	UTM N:	b.	UTM E:				
2. Check which, if any, of th	ne Tier I inclusion	ary criteria are m	net by the Dispo	osal Site, pursua	ant to 310 CMR 40.05	520(2):	
a. Groundwater is loc contamination by an C the applicable RCGW	Oil or Hazardous /-1 Reportable Co	Material at the ting the modern of the moder	me of Tier Clas forth in 310 CM	sification at cor			
C. EXPOSURE PATHWAYS	•						
 Exposure Pathways, and For A. through D., score ac 			•	av Designation	Criteria and NRS Tah	ale II	
For E., score using NRS To		WIIX 40.1012 EA	cposure r diriwi	ay Designation		710 II.	
				Score	1		
	A. Soil (include	es sediment)					
	B. Groundwate	er					
	C. Surface Wa	ter (includes we	etlands)				
D. Air							
	E. Number of C	OHM Sources					
	Total NRS Sec	tion II Score (15	i - 700)				
2. Was Section G (NRS Se	ection VI) used to	amend the score	e for this Section	n of the NRS?	a. Yes [b. No	

Revised: 02/10/2004 Page 1 of 5



NUMERICAL RANKING SYSTEM (NRS) SCORESHEET

BWSC	107A
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Release	Tracking Number
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7/15	Pursuant to 310 CMR 40.1511 (Subpart O)						
3. Summary Rationale for Exposure Pathway Values, A. through D., and Phase I Report References:							
	CHARACTERISTICS (NRS SECTION II us Material (OHM) Toxicity Score (NRS	-	ection III.A.):				
a. List the Four	Highest OHM Toxicity Scores from NF	RS	Table III.A.:				
	OHM Scored		Concentratio	n and Media		Toxicity Sc	ore (1 - 80)
b. Score using	NRS Worksheet III.A.1. to determine the	ne	OHM Toxicity Score for	or OHM not listed ir	NRS	S Table III.A.:	
	ОНМ	Н	uman Health-based Toxicity Value	Concentration (Soil - ug/g)		oncentration Vater - ug/l)	Toxicity Score
c. Use the Hig	hest OHM Toxicity Score from eithe		NRS Table III.A. or \	1		7	
	OHM Scored			Toxicity Scor	е	1	

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Release	Tracking Number

Pursuant to 310 C			•	K5) 50	ORE	SHEET				
2. Multiple OHMs (NRS Section III.B.):										
Was the Toxicity Score of more than one OHM greater than or equal to 30? a. Yes (30) b. No (0)										
3. OHM Mobility and Persistence (NRS Section III.C.):										
Score according to 310 CMR 40.	1514 - OHM	Mobilit	ty and Persis	tence						
a. OHM So	cored			b. Score (0 - 50)						
4. Disposal Site Hydrogeology (NRS S	Section III.D.):		Site Hydr		gy				
Score according to 310 CMR 40.1515 - Soil Permeability, and NRS Table III.D.										
5. Total NRS Section III Score:	Α.	В	. C.	ı	Э.		or Section II 3 - 180)	I]	
6. Was Section G (NRS Section VI) u	sed to amen	d the s	score for this	Section o	f the N	RS?	a. Yes		b. No	
E. HUMAN POPULATION AND LAND	USES (NRS \$	SECTIO	ON IV):							
Human Population (NRS Section IV Score using NRS Table IV.A.	.A.):		ential Popula thin 1/2 Mile	tion		utions 500 Feet	On-site Workers		opulation ore (0 - 40)	
2. Aquifers (NRS Section IV.B.):										
a. Sole Source Aquifer:	i. Yes (25	i) ii.	Name:					iii. No	o (0)	
b. Potentially Productive Aquifer:		i. Med	ium or High (15)	ii	i. No (0)				
3. Water Use (NRS Section IV.C.):	Proximity		Persons Se			ate Water	Alternate Pu		Water	
Score using NRS Table IV.C.	Public Drin Water Sou		Public Dr Water S			upplies n 500 Feet	Water Supp Available		Use Score (0 -125)	
4. Total NRS Section IV Score:	A.		B.	C.			r Section IV - 205)			
5. Was Section G (NRS Section VI) u	sed to amen	d the s	score for this	Section o	f the N	RS?	a. Yes [t	o. No	

Page 3 of 5 Revised: 02/10/2004



BW	SC	10	7A
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			AL RANKING \$ 10 CMR 40.1511 (Su	SYSTEM (NRS) abpart O)	SCO	ORESHEE	ET '	Release	e Tra -	cking Number	
F.	ECOLOGICAL PO	OPULATION (N	RS SECTION V):								
1.	Environmental Re	esource Areas	(NRS Section V.A.):	Score using NRS	Table	V.A.					
	Area of Critical Environmental Concern	Threatened	Special Concern, d or Endangered ies Habitat	Wetlands, Certified Pool, or Outstan Resource Wat	ding	Fish Habitat	Protected Open Space		Res	ironmental ource Area re (0 - 150)	
2.	Environmental To	xicity Score (N	RS Section V.B.):								
	Score only if En	vironmental R	esource Area Score	is greater than or ed	ual to	30.					
	a. List the Thre	e Highest Env	ironmental Toxicity S	Scores from NRS Tab	le V.E	3.:					
		ОН	IM Scored		C	Concentration	n and Med	dia	T	oxicity Score (0 - 35)	
				e the Environmental /alues for each OHM		ty Score for C	OHM not li	isted in	NRS	Table V.B.	
		ОНМ		Environmer Toxicity Val		Concentration (Soil - ug/g		centrat iter - ug		Environmenta Toxicity Score	
	c. Use the Highe	est Environme	ntal Toxicity Score f	om either NRS Table	V.B.	or from Work	sheet V.B	3.1.:			
	Г		OHM Sco	ored		Toxicit	ty Score	1			
								•			
3.	Total NRS Section	n V Score:	A.	B.	Tota	I for Section (0 - 185)	n V				
4.	Was Section G (I	NRS Section V	/I) used to amend th	e score for this Section	on of t	he NRS?	a. Ye	s 🔲	b.	No	

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Release Tracking Number

BWSC107A

NUMERICAL RANKING SYSTEM (NRS) SCORESHEET Pursuant to 310 CMR 40.1511 (Subpart O)

G. MITIGATING DISPOSAL SITE -SPECIFIC CONDITIONS (NRS SECTION VI):
1. Disposal site-specific conditions that warrant amending the site score. Changes directly related to NRS Sections or Subsection scores may not reduce the score more than the relevant subsection value assigned for the disposal site in that subsection. Section VI must reference specific pages of the Phase I. Section VI may not exceed plus or minus 50 points and may be scored only in 5-point increments.
2. Check here if additional pages are provided in an attachment.
3. Disposal Site Amendment (Not to exceed plus or minus 50 points): Total Score Section VI

Revised: 02/10/2004 Page 5 of 5



Massachusetts Department of Environmental Protection

Bureau of Waste Site Cleanup TIER II COMPLIANCE HISTORY Pursuant to 310 CMR 40.0540 (Subpart E)			Release Tracking Number
A. DISPOSAL SITE COMPLIANCE HISTORY SUN	IMARY:		
Check here if a Tier II Compliance Histo there has been no change in that person's completed.			
2. List all permits or licenses that have been is	sued by the Departme	nt that are relevant to this Disp	oosal Site:
Program	Permit Number	Permit Category	Facility ID
a. Air Quality			
b. Hazardous Waste (M.G.L. c. 21C)			
c. Solid Waste			
d. Industrial Wastewater Management			
e. Water Supply			
f. Water Pollution Control/Surface Water			
g. Water Pollution Control/Groundwater			
h. Water Pollution Control/Sewer Connection			
i. Wetland & Waterways			
List all other Federal, state or local permits, li this Disposal Site:	censes, certifications,	registrations, variances, or app	provals that are relevant to
Issuing Authority or Program, or Documentation Type		Identification Number	Date Issued mm/dd/yyyy
			_
4. Check here to certify that, if needed, a sattached. This statement must describe the compliance history or regulations; and (2) other laws for the protection of he government agency. Such a statement should identify enforce its requirements including, but not limited to, (PAN), Notice of Intent to Take Response Action (NOR).	of the person or entity na ealth, safety, public welfa / information such as: (1 a Notice of Noncomplian A), and an administrative	med in BWSC107, Section D with a re and the environment administer) actions relevant to the Disposal ce (NON), Notice of Intent to Asse enforcement order; (2) administr	the following: (1) DEP red or enforced by any other Site taken by the Department to ess Civil Administrative Penalty ative consent orders; (3) judicial
consent judgements; (4) similar administrative action the Disposal Site brought on behalf of the DEP or other action identified provide the following information: (4)	er Federal, state, or local	agencies; and (6) any additional re	elevant information. For each

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description of noncompliance cited; (3) current status of the matter; and (4) final disposition, if any.



APPENDIX F

Public Notification Documentation



May 14, 2008.

Section Chief Bureau of Waste Site Cleanup Massachusetts Department of Environmental Protection Central Region Main Office 627 Main Street Worcester, MA 01608

RE: Copy of Legal Notice

Tier Classification and Phase I - Initial Site Investigation Report

Former Nu-Style Company, Inc.

87 Grove Street

Franklin, Massachusetts

RTN: 2-16694

Dear Section Chief:

In accordance with the Massachusetts Contingency Plan (MCP; 310 CMR 40.1403 (6)), Fuss & O'Neill, Inc. has prepared this letter on behalf of the Town of Franklin, Massachusetts, the owner of the subject property, to provide you with a copy of the legal notice published in The Milford Daily News, Milford, Massachusetts on May 19, 2008. The legal notice indicated the Tier Classification for the release identified by the Massachusetts Department of Environmental Protection (MADEP) as Release Tracking Number (RTN) 2-16694. The site has been classified as a Tier II. A copy of the legal notice is attached. Copies of the cover letter sent to the Chief Municipal Officer and Health Agent of the City of Attleboro are also attached.

Sincerely

317 Iron Horse Way Suite 204

Providence, RI 02908 David I.P. Foss

Senior Project Manager

t (401) 861-3070 (800) 286-2469

f (401) 861-3076

Attachment: Legal Notice

www.FandO.com Cover Letter to Town Administrator Jeffrey Nutting Cover Letter to Health Agent David E. McKearney, R.S.

Connecticut

Massachusetts

Town of Franklin

New York Rhode Island

Ms. Stephanie Mercandetti, Norfolk County

North Carolina

South Carolina

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Liz Cassidy Franklin Planning Board 355 East Central Street Franklin, Ma 02038

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Milford Daily News

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5/19/2008

Stop date: #Lines:

5/19/2008

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\$93.15

Insertions:

Pmt. Type:

Invoice Total: \$93.15

Payment Information

Receipt#

87 Grove Street

LEGAL NOTICE NOTICE OF INITIAL SITE **INVESTIGATION AND** TIER II CLASSIFICATION

FORMER NU-STYLE PROPERTY 87 GROVE STREET, FRANKLIN, MASSACHUSETTS **RELEASE TRACKING NUMBER 2-16694**

A release of oil and/or hazardous materials has occurred at this location, which is a disposal site as defined by M.G.L. c. 21E, § 2 and the Massachusetts Contingency Plan, 310 CMR 40.0000. To evaluate the release, a Phase I Initial Site Investigation was performed pursuant to 310 CMR 40.0480. As a result of this investigation, the site has been classified as Tier II pursuant to 310 CMR 40.0500. On May 14, 2008 the Town of Franklin, Massachusetts filed a Tier II Classification Submittal with the Department of Environmental Protection (MassDEP). To obtain more information on this disposal site, please contact Mr. David Foss, Fuss & O'Neill, Inc., 317 Iron Horse Way, Suite 204, Providence, RI 02908, 401-861-3070 ext. 4579.

The Tier II Classification Submittal and the disposal site file can be reviewed at MassDEP Central Region Main Office, 627 Main Street, Worcester, MA 01608, 508-792-7650.

Additional public involvement opportunities are available under 310 CMR 40.1403(9) and 310 CMR 40.1404.

AD#11669992 MDN 5/19/08



May 14, 2008

Jeffrey Nutting Town Administrator Franklin Town Hall 355 East Central Street Franklin, MA 02038

Re:

Notice of Availability of Tier Classification and Phase I - Initial Site Investigation Report and

Copy of Legal Notice

Former Nu-Style Company, Inc.

87 Grove Street

Franklin, Massachusetts

RTN: 2-16694

Dear Mr. Nutting:

Fuss & O'Neill, Inc. has prepared this letter on behalf of the Town of Franklin, Massachusetts, the owner of the subject property, to provide notice in accordance with the Massachusetts Contingency Plan (MCP; 310 CMR 40.0000) that a Tier Classification and Phase I – Initial Site Investigation Report has been submitted to the Massachusetts Department of Environmental Protection (MADEP) for the release identified by the MADEP as Release Tracking Number (RTN) 2-16694. The documents indicate that volatile organic compounds and metals were identified in on-site soil and groundwater and polycyclic aromatic hydrocarbons were identified in on-site sediment. The site has been classified as a Tier II.

The reports and supporting documentation are available for review or to be copied at the following location:

Massachusetts Department of Environmental Protection Central Region Main Office 627 Main Street Worcester, MA 01608

Telephone: 508-792-7650

317 Iron Horse Way Suite 204

Providence, RI 02908

In accordance with the MCP (310 CMR 40.1403 (6)), Fuss & O'Neill, Inc. has also prepared this letter to provide you with a copy of the legal notice published in The Milford Daily News, Milford, Massachusetts on May 19, 2008. The legal notice indicated the Tier Classification for the release identified above. A copy of the legal notice is attached.

t (401) 861-3070 Sincerely

(800) 286-2469 f (401) 861-3076

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86-2469

David J.P. Foss Senior Project Manager

www.FandO.com

MADEP

Connecticut

Attachment:

Legal Notice

Massachusetts

New York

Rhode Island

:

Ms. Stephanie Mercandetti, Norfolk County

North Carolina

South Carolina

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May 14, 2008

David E. McKearney, R.S. Health Agent/Director Franklin Town Hall 355 East Central Street Franklin, MA 02038

Re:

Notice of Availability of Tier Classification and Phase I - Initial Site Investigation Report and

Copy of Legal Notice

Former Nu-Style Company, Inc.

87 Grove Street

Franklin, Massachusetts

RTN: 2-16694

Dear Mr. McKearney:

Fuss & O'Neill, Inc. has prepared this letter on behalf of the Town of Franklin, Massachusetts, the owner of the subject property, to provide notice in accordance with the Massachusetts Contingency Plan (MCP; 310 CMR 40.0000) that a Tier Classification and Phase I – Initial Site Investigation Report has been submitted to the Massachusetts Department of Environmental Protection (MADEP) for the release identified by the MADEP as Release Tracking Number (RTN) 2-16694. The documents indicate that volatile organic compounds and metals were identified in on-site soil and groundwater and polycyclic aromatic hydrocarbons were identified in on-site sediment. The site has been classified as a Tier II.

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Telephone: 508-792-7650

In accordance with the MCP (310 CMR 40.1403 (6)), Fuss & O'Neill, Inc. has also prepared this letter to provide you with a copy of the legal notice published in The Milford Daily News, Milford, Massachusetts on May 19, 2008. The legal notice indicated the Tier Classification for the release identified above. A copy of the legal notice is attached.

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Sincerely,

David J.P. Foss

Senior Project Manager

MADEP

Connecticut

ticut Attachment:

Legal Notice

Massachusetts

New York

C:

Ms. Stephanie Mercandetti, Norfolk County

Rhode Island

North Carolina

South Carolina

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