



Wilcox & Barton INC.

ENVIRONMENTAL AND ENGINEERING SERVICES

**BIENNIAL GROUNDWATER MONITORING
REPORT, 2011**

**FORMER WAIT FARM MOTOR INN
4805 MAIN STREET
WAITSFIELD, VERMONT
VT DEC SITE NUMBER 2002-2982**

Prepared for:

Gumbo Group, LLC
4805 Main Street
Waitsfield, Vermont 05673
Contact: Mr. Branwell Lepp, (802) 583-5525

Prepared by:

Wilcox & Barton, Inc.
57 Hoit Road
Concord, New Hampshire 03301
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September 2, 2011

Wilcox & Barton, Inc. Project No.: GUMB0001

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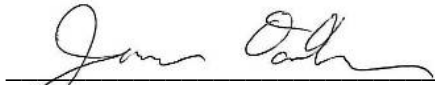


CERTIFICATION

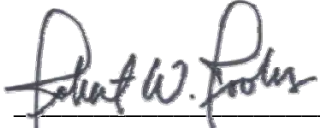
The following personnel have prepared and/or reviewed this report for accuracy, content, and quality of presentation.

Document Title: Biennial Groundwater Monitoring Report, 2011
Former Wait Farm Motor Inn
4805 Main Street, Waitsfield, Vermont
VT DEC Site #2002-2982

Date/Version: September 2, 2011



James Donaldson
Project Scientist



Robert W. Rooks, PE
Principal Engineer

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1.0 INTRODUCTION

This report presents the results of a biennial groundwater sampling event performed at the former Wait Farm Motor Inn, 4805 Main Street in Waitsfield, Vermont (the site). The groundwater monitoring activities are a continuation of response actions implemented following the discovery of a release of an unknown quantity of #2 fuel oil from an underground storage tank (UST) supply line. The leak was discovered following UST system removal in March 2002. The Vermont Department of Environmental Conservation (VT DEC) assigned Site Number 2002-2982 to the release.

The property was sold in March 2010 and Wilcox & Barton, Inc. was retained by Mr. Branwell Lepp of Gumbo Group, LLC, in May 2011 to provide environmental consulting services. The current biennial groundwater monitoring event was performed in accordance with the VT DEC letter dated October 9, 2009, and a scope of work and budget approved by VT DEC in a letter dated June 3, 2011. A copy of the October 9, 2009, VT DEC correspondence is included in Appendix A.

The location of the site is presented on Figure 1 – *Site Location Map*, and pertinent site features are depicted on Figure 2 – *Site Plan*. A brief history of response actions at the site to date is provided in Appendix B.

2.0 SITE DESCRIPTION

The site is located in a rural area of Waitsfield, Vermont, consisting primarily of commercial properties situated along Main Street (Vermont Route 100). The site is operated as a retail showroom and office space for a residential and commercial heating, ventilation, and air conditioning supply company. The site is occupied by a two-story, wood-frame structure containing apartments, a showroom, and office space in the full basement. Unimproved portions of the property consist of unpaved parking areas, lawn, and landscaping. A pond is located on the eastern portion of the property.

The topography of the western portion of the property between the building and Main Street is predominantly level. The site slopes steeply downward east of the building toward the pond. The site is serviced by an on-site septic system and private drinking water well. The leach field for the septic system is located between the building and Main Street, and is shared with the adjacent residential property. According to prior site investigations, the drinking water well is 300 feet deep, is installed in bedrock, and is located approximately 250 feet northeast of the release area. The water supply well is also shared with the adjacent residential property.

3.0 RESPONSE ACTIONS

Response actions conducted during this reporting period consist of a groundwater monitoring event performed on June 16, 2011. A summary of monitoring activities is presented in this section.

3.1 Groundwater Gauging

On June 16, 2011, monitoring wells MW-1, MW-2, and MW-4 were gauged for depth to water using an oil/water interface probe capable of measuring depth to water and non-aqueous phase liquid (NAPL) to the nearest 0.01 foot. Monitoring well MW-3 could not be located on the densely vegetated slope, which showed signs of disruption from the conversion of the building from an Inn to office space. The well is presumed to have been destroyed.

The presence of NAPL was investigated by inserting a clean, clear bailer into each well until half the volume of the bailer was filled, removing the bailer, and visually inspecting the contents. Gauging did not reveal a measurable thickness of NAPL in any well. Well gauging data and product thickness measurements are summarized in Table 1.

A petroleum sheen was observed on the purge water from well MW-2, and a petroleum odor was noted in purge water from wells MW-2 and MW-4 prior to sampling. The well riser for MW-2 was damaged and would not accept a probe or bailer, so the riser was cut back to facilitate well gauging and sampling.

3.2 Groundwater Sample Collection and Analysis

On June 16, 2011, groundwater samples were collected from monitoring wells MW-1, MW-2, and MW-4. Sampling for volatile organic compounds (VOCs) was performed in accordance with Wilcox & Barton, Inc. Field Sampling Protocol, a copy of which is included in Appendix C. The samples were submitted under Chain of Custody to Con-Test Analytical Laboratory of East Longmeadow, Massachusetts, for analysis of VOCs by U.S. Environmental Protection Agency (EPA) Method 8021B.

One or more target constituents were detected in monitoring wells MW-1, MW-2, and MW-4. The samples from MW-1, MW-2, and MW-4 were found to contain several VOCs indicative of fuel oil, although none were present at concentrations above Vermont Primary Groundwater Enforcement Standards (VGES). Detected concentrations of fuel oil-related constituents generally show a decreasing concentration trend from previous monitoring results.

Four constituents were detected for the first time in any site well during this monitoring event: acetone, tert-butylbenzene, chloromethane, and trichlorofluoromethane. All detected concentrations of acetone and trichlorofluoromethane were below the corresponding VGES; no VGES is available for tert-butylbenzene. The detected concentrations of chloromethane exceeded the VGES of 30 micrograms per liter ($\mu\text{g/L}$) in all three wells, at concentrations ranging from 37 $\mu\text{g/L}$ in MW-4 to 110 $\mu\text{g/L}$ in MW-2. The source of the chloromethane is currently unknown.

Graphs of concentration trends for Total VOCs and naphthalene are presented in Appendix D. Stable to downward trends are evident over the two year period from 2009 to 2011. Only MW-2 shows a slight upwards trend in Total VOCs; however, the concentrations are well below those found during the historical high period in 2002 and 2003 and are largely attributable to the presence of chloromethane. The concentration of naphthalene in MW-2 remains stable.

Dissolved phase constituent concentrations across the remainder of site have shown a decrease from historic highs, and NAPL has not been detected since 2008.

A summary of the groundwater data is presented in Table 2 and a copy of the laboratory report is included in Appendix E.

4.0 POTENTIAL SENSITIVE RECEPTORS AND EXPOSURE PATHWAYS

In general, potential human receptors may include residents, workers, visitors, or trespassers in the vicinity of the site. Environmental receptors may include flora and fauna within the affected area. Potential exposure pathways and points may include wellhead protection areas, potable water wells, surface waters, wetlands, migration of vapors into building through basements, direct contact (e.g., playgrounds), and utility corridors. Exposure routes through which contamination may reach human receptors include ingestion, inhalation, and dermal contact.

Based on a review of response actions undertaken by Griffin and a limited inspection of the property, Wilcox & Barton, Inc. has identified the following potentially complete exposure pathways for human receptors at the site:

Ingestion of Drinking Water - The site and surrounding properties are serviced by private drinking water wells. No municipal water service is provided within the Town of Waitsfield. As groundwater contamination has been documented at the site, drinking water from wells located within the area of contamination may serve as a potential exposure point. The potable well located on the subject property is reportedly located approximately 250 feet northeast of (cross-gradient from) the source area and is installed in bedrock at a depth of 300 feet. The potable well was sampled and analyzed for VOCs by EPA Method 524.2 on June 20, 2002. The analysis did not detect fuel oil constituents.

According to the Griffin International “*Initial Report on the Investigation of Subsurface Petroleum Contamination and Product Recovery Efforts*” report of August 28, 2002, no known private or public water supply wells, other than the well on the property, are located within 500 feet downgradient of the site.

Inhalation of Vapors – The former Wait Farm Motor Inn building is constructed over a basement, which is now occupied by employees during working hours. The release of fuel oil occurred beneath the building. NAPL has been detected within shallow groundwater (3 to 4 feet below the basement floor), most recently during the 2008 sampling event. In addition, in an inspection of the basement during the UST closure in March 2002, screening of the indoor air with a photoionization detector (PID) identified a peak VOC concentration of 2.8 parts per million by volume (ppmv). Constituent concentrations in soil and groundwater and VOC screening measurements within the basement suggest that impacts to indoor air may exist, although the influence of two 275-gallon above-ground storage tanks and an oil-fired boiler in the basement cannot be ruled out.

Dermal Contact with Soil, Groundwater, Surface Water – Based on historic site characterization data, contamination is present in soil and groundwater beneath the basement floor and in soil and

groundwater within the vicinity of the water table extending to the on-site pond. The pond was historically impacted with product, however, no sheen was visible on the pond surface during the 2011 monitoring event. Contact with surface water may result in dermal exposures. Under current conditions, dermal contact with soil and groundwater are not likely unless excavations are performed.

Impact to Surface Water - The pond on the subject property has historically been impacted with petroleum product. Product recovery efforts were instituted on the pond using absorbent pads and containment booms. Due to the concentration of constituents remaining within groundwater and the location of the pond relative to the source area, dissolved-phase constituents may be impacting the pond. No surface water samples have been collected to document the potential impacts.

5.0 CONCLUSIONS AND RECOMMENDATIONS

During this reporting period, a biennial groundwater monitoring event was performed on July 16, 2011. Gauging and purging of monitoring wells did not identify the presence of NAPL, but a petroleum sheen was identified on purge water from well MW-2. VOCs consistent with fuel oil were measured in wells MW-1, MW-2, and MW-4, but no fuel oil-related constituents were detected at concentrations exceeding VGES. Chloromethane was detected in samples from wells MW-1, MW-2, and MW-4 for the first time since groundwater monitoring began at the site, and all detected concentrations exceeded the VGES for chloromethane. The source of chloromethane is currently unknown.

Although the current monitoring data did not indicate the presence of NAPL in the subsurface, historic data suggest that it is still present in some quantity. NAPL represents a continuing source of both dissolved and vapor phase contamination. Such conditions pose a threat of impact to drinking water, surface water, and indoor air at the site. Drinking water quality data have not been collected since 2002.

Wilcox & Barton, Inc. recommends the continuation of groundwater and NAPL gauging, groundwater monitoring for VOCs, and reporting on a biennial basis. Due to the historical presence of contaminant concentrations exceeding VGES in monitoring well MW-3, Wilcox & Barton, Inc. recommends that the monitoring well be replaced prior to the scheduled 2013 monitoring event. The well riser for MW-2 should also be repaired.

TABLES



TABLE 1
Well Gauging and Piezometric Head Elevation Data
 Wait Farm Motor Inn
 4805 Main Street, Waitsfield, Vermont

| Well ID | Gauging Date | Depth to Water (ft) | Depth to NAPL (ft) | NAPL Thickness (ft) | Well Elevation (Ft) | Piezometric Head Elevation (ft) |
|----------|--------------|---------------------|--------------------|---------------------|---------------------|---------------------------------|
| MW-1 | 11/21/02 | 3.28 | 3.16 | 0.125* | NM | NM |
| | 02/10/03 | 3.45 | -- | -- | NM | NM |
| | 05/21/03 | 3.50 | -- | -- | NM | NM |
| | 09/25/03 | 3.41 | -- | -- | NM | NM |
| | 03/29/04 | 3.38 | -- | -- | NM | NM |
| | 05/27/05 | 3.28 | 3.26 | 0.02 | NM | NM |
| | 03/29/06 | 4.50 | 4.48 | 0.02 | NM | NM |
| | 07/21/06 | 3.02 | 3.00 | 0.02 | NM | NM |
| | 10/21/06 | 2.93 | 2.92 | 0.01 | NM | NM |
| | 01/24/07 | 3.09 | -- | -- | NM | NM |
| | 05/02/07 | 7.55 | -- | -- | NM | NM |
| | 05/06/08 | 3.61 | -- | -- | NM | NM |
| | 05/18/09 | 3.15 | -- | -- | NM | NM |
| | 06/16/11 | 3.53 | -- | -- | NM | NM |
| | MW-2 | 11/21/02 | 1.48 | -- | -- | NM |
| 02/10/03 | | 2.36 | -- | -- | NM | NM |
| 05/21/03 | | 2.55 | -- | -- | NM | NM |
| 09/25/03 | | 2.21 | -- | -- | NM | NM |
| 03/29/04 | | 1.88 | -- | -- | NM | NM |
| 05/27/05 | | 2.31 | -- | -- | NM | NM |
| 03/29/06 | | 2.10 | -- | -- | NM | NM |
| 07/21/06 | | 2.18 | -- | -- | NM | NM |
| 10/21/06 | | 2.88 | -- | -- | NM | NM |
| 01/24/07 | | 1.90 | -- | -- | NM | NM |
| 05/02/07 | | 1.83 | -- | -- | NM | NM |
| 05/06/08 | | 2.20 | -- | -- | NM | NM |
| 05/18/09 | | 1.77 | -- | -- | NM | NM |
| 06/16/11 | | 0.99 | -- | -- | NM | NM |
| MW-3 | | 11/21/02 | 3.76 | 3.64 | 0.125* | NM |
| | 02/10/03 | 4.98 | -- | -- | NM | NM |
| | 05/21/03 | 4.28 | -- | -- | NM | NM |
| | 09/25/03 | 4.55 | -- | -- | NM | NM |
| | 03/29/04 | 2.30 | -- | -- | NM | NM |
| | 05/27/05 | 4.52 | -- | -- | NM | NM |
| | 03/29/06 | 4.69 | -- | -- | NM | NM |
| | 07/21/06 | 3.96 | -- | -- | NM | NM |
| | 10/21/06 | 3.81 | -- | -- | NM | NM |
| | 01/24/07 | 4.50 | -- | -- | NM | NM |
| | 05/02/07 | 3.98 | 3.97 | 0.01 | NM | NM |
| | 05/06/08 | 4.68 | -- | -- | NM | NM |
| | 05/18/09 | 3.32 | -- | -- | NM | NM |
| | 06/16/11 | | | | Well not found | |
| | MW-4 | 11/21/02 | 6.55 | -- | -- | NM |
| 02/10/03 | | 6.97 | -- | -- | NM | NM |
| 05/21/03 | | 6.67 | -- | -- | NM | NM |
| 09/25/03 | | 6.55 | -- | -- | NM | NM |
| 03/29/04 | | 6.54 | -- | -- | NM | NM |
| 05/27/05 | | 6.50 | -- | -- | NM | NM |
| 03/29/06 | | 6.75 | -- | -- | NM | NM |
| 07/21/06 | | 6.38 | -- | -- | NM | NM |
| 10/21/06 | | 6.20 | -- | -- | NM | NM |
| 01/24/07 | | 6.54 | -- | -- | NM | NM |
| 05/02/07 | | 4.55 | -- | -- | NM | NM |
| 05/06/08 | | 6.64 | -- | -- | NM | NM |
| 05/18/09 | | 6.21 | -- | -- | NM | NM |
| 06/16/11 | | 7.71 | -- | -- | NM | NM |

Notes:

Data prior to 5/27/05 collected by Griffin International, Inc.

* 1/8 inch of NAPL measured using a bailer by Griffin International, Inc.

NAPL Non-aqueous phase liquid.

NM Not Measured.



TABLE 2
Groundwater Samples - Summary of Analytical Results
 Wait Farm Motor Inn
 4805 Main Street, Waitsfield, Vermont

| Sample Identification Sample Date | Vermont Groundwater Enforcement Standard † | MW-1 | | | | | | | | | | | |
|--|---|------------|----------|------------|-------------|-------------|-------------|----------|----------|-----------------|----------|----------|-----------|
| | | 07/03/02 | 11/21/02 | 02/10/03 | 05/21/03 | 09/25/03 | 03/29/04 | 05/27/05 | 03/29/06 | 05/02/07 | 05/07/08 | 05/18/09 | 06/16/11 |
| Volatile Organic Compounds (VOCs) by EPA Method 8021B/8260B | | | | | | | | | | | | | |
| Acetone | 700 | / | / | / | / | / | / | NS* | NS* | < 500 | NS* | < 50 | 50 UB |
| tertiary-Amyl methyl ether (TAME) | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | <5.0 | NS* | <0.50 | 0.50 U |
| Benzene | 5.0 | 5.9 | NS* | ND | ND | ND | ND | NS* | NS* | <10.0 | NS* | <1.0 | 1.0 U |
| n-Butylbenzene | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | 16.4 | NS* | <1.0 | 1.0 U |
| sec-Butylbenzene | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | 11.8 | NS* | <1.0 | 0.14 J |
| tert-Butylbenzene | NE | / | / | / | / | / | / | NS* | NS* | < 10.0 | NS* | < 1.0 | 1.0 U |
| Chloromethane | 30.0 | / | / | / | / | / | / | NS* | NS* | < 20.0 | NS* | < 2.0 | 38 |
| Ethylbenzene | 700.0 | 41.3 | NS* | ND | ND | ND | ND | NS* | NS* | 113 | NS* | <1.0 | 1.0 U |
| Isopropylbenzene | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | 29.4 | NS* | <1.0 | 1.0 U |
| p-Isopropyltoluene | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | <10.0 | NS* | <1.0 | 1.0 U |
| Methyl tert-butyl ether (MTBE) | 40.0 | 31.8 | NS* | ND | ND | ND | ND | NS* | NS* | 12.5 | NS* | <1.0 | 0.29 J |
| Naphthalene | 20.0 | 127 | NS* | 121 | 63.6 | 54.1 | 96.8 | NS* | NS* | 30.5 | NS* | <2.0 J | 2.0 UJ |
| n-Propylbenzene | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | 48.8 | NS* | <1.0 | 1.0 U |
| 1,1,2,2-Tetrachloroethane | NE | -- | NS* | -- | -- | -- | -- | NS* | NS* | <5.0 | NS* | <0.50 | 0.50 U |
| Toluene | 1,000.0 | 42.8 | NS* | ND | ND | ND | ND | NS* | NS* | <10.0 | NS* | <1.0 | 1.0 U |
| Trichlorofluoromethane | 2,100.0 | / | / | / | / | / | / | NS* | NS* | < 20.0 | NS* | < 2.0 | 1.3 J |
| 1,2,4-Trimethylbenzene | NE | 235 | NS* | 49.8 | 20.2 | 39.8 | 81.3 | NS* | NS* | 176 | NS* | <1.0 | 1.0 U |
| 1,3,5-Trimethylbenzene | NE | 70.6 | NS* | 25.7 | ND | 12.7 | 25.5 | NS* | NS* | 83.1 | NS* | <1.0 | 1.0 U |
| Total Trimethylbenzenes | 350 | 249 | NS* | 121.0 | 63.6 | 54.1 | 96.8 | NS* | NS* | 262.4 | NS* | <2.0 | 2.0 U |
| Xylenes (total) | 10,000 | 298 | NS* | 19.9 | ND | ND | 14.8 | NS* | NS* | 169.3 | NS* | <3.0 | 3.0 U |
| Total Petroleum Hydrocarbons (TPH) by EPA Method 8015 | | | | | | | | | | | | | |
| TPH | NE | -- | NS* | 12,300 | 16,100 | 25,500 | 147,000 | NS* | NS* | -- | NS* | -- | -- |

All results are in micrograms per liter (ug/L).

Only detected and selected other analytes shown; all others not detected.

- bold shaded** Exceeds Vermont Groundwater Enforcement Standard (VGES)
- bold italics** Reporting limit for non-detection exceeds VGES.
- "--" Not analyzed.
- ND Not detected; reporting limit not shown.
- <X; U Not detected; reporting limit shown.
- NS* Not sampled, free product present.
- NE No standard established.
- J Estimated concentration below reporting limit or due to quality control limitations.
- B Amount in sample >5x (>10x for acetone) amount in method blank; result valid.
- UB Amount in sample <5x (<10x for acetone) amount in method blank; changed to non-detection.
- "/" Analytical result not available.
- † Groundwater Protection Rule and Strategy, 02/05, and Interim Groundwater Quality Standards, Guidance, 08/07.



TABLE 2
Groundwater Samples - Summary of Analytical Results
 Wait Farm Motor Inn
 4805 Main Street, Waitsfield, Vermont

| Sample Identification Sample Date | Vermont Groundwater Enforcement Standard † | MW-2 | | | | | | | | | | | |
|--|---|--------------|------------|-------------|----------|-------------|-------------|------------|------------|----------|----------|----------|------------|
| | | 07/03/02 | 11/21/02 | 02/10/03 | 05/21/03 | 09/25/03 | 03/29/04 | 05/27/05 | 03/29/06 | 05/02/07 | 05/07/08 | 05/18/09 | 06/16/11 |
| Volatile Organic Compounds (VOCs) by EPA Method 8021B/8260B | | | | | | | | | | | | | |
| Acetone | 700 | / | / | / | -- | / | / | < 10.0 | < 50.0 | < 50.0 | < 50 | < 50 | 50 UB |
| tertiary-Amyl methyl ether (TAME) | NE | -- | -- | -- | -- | -- | -- | 1.0 | 0.5 J | <0.5 | <0.5 | <0.50 | 0.50 U |
| Benzene | 5.0 | 20.3 | ND | ND | -- | ND | 11.9 | 7.2 | 5.8 | 3.6 | 1.4 | 1.5 | 1.2 |
| n-Butylbenzene | NE | -- | -- | -- | -- | -- | -- | 2.6 | <1.0 | 1.2 | 1.3 | <1.0 | 2.1 |
| sec-Butylbenzene | NE | -- | -- | -- | -- | -- | -- | < 0.6 | <1.0 | 1.9 | 1.8 | 1.2 | 4.5 |
| tert-Butylbenzene | NE | / | / | / | -- | / | / | < 0.8 | < 1.0 | < 1.0 | < 1.0 | < 1.0 | 1.0 U |
| Chloromethane | 30.0 | / | / | / | -- | / | / | < 1.2 | < 2.0 | < 2.0 | < 2.0 | < 2.0 | 110 |
| Ethylbenzene | 700.0 | 11.2 | 16.3 | ND | -- | ND | ND | 12.3 | 28.1 | 30.9 | 21.4 | 7.3 | 7.0 |
| Isopropylbenzene | NE | -- | -- | -- | -- | -- | -- | 2.5 | 5.4 | 8.3 | 8.3 | 5.0 | 9.9 |
| p-Isopropyltoluene | NE | -- | -- | -- | -- | -- | -- | 0.9 | <1.0 | <1.0 | <1.0 | <1.0 | 0.41 J |
| Methyl tert-butyl ether (MTBE) | 40.0 | 353 | 111 | 61.0 | -- | 30.7 | 33.6 | 16.7 | 8.9 | 3.9 | 1.8 | 2.1 | 1.3 |
| Naphthalene | 20.0 | 48.7 | 131 | 20.5 | -- | 49.2 | ND | 3.6 | 6.8 J | 10.5 | 8.7 | 4.7 | 5.0 JB |
| n-Propylbenzene | NE | -- | -- | -- | -- | -- | -- | 1.6 | 5.1 | 8.7 | 8.5 | 3.9 | 8.9 |
| 1,1,2,2-Tetrachloroethane | NE | -- | -- | -- | -- | -- | -- | < 0.5 | <0.5 | 0.8 | <0.5 | <0.50 | 0.50 U |
| Toluene | 1,000.0 | 4.2 | ND | ND | -- | ND | ND | < 0.7 | <1.0 | <1.0 | <1.0 | <1.0 | 0.34 J |
| Trichlorofluoromethane | 2,100.0 | / | / | / | -- | / | / | < 0.7 | < 2.0 | < 2.0 | < 2.0 | < 2.0 | 2.0 U |
| 1,2,4-Trimethylbenzene | NE | 54.9 | 114 | 44.2 | -- | 11.1 | ND | 0.8 | 3.1 | 12.4 | 11.6 | 2.4 | 6.1 |
| 1,3,5-Trimethylbenzene | NE | 29.4 | 84.7 | 114 | -- | 144 | 11.3 | 2.6 | 1.9 | 3.4 | 5.8 | <1.0 | 1.0 U |
| Total Trimethylbenzenes | 350 | 437.4 | 258 | 81.5 | -- | 79.9 | 45.5 | 48.4 | 53.3 | 69.8 | 53.2 | 2.4 | 6.1 |
| Xylenes (total) | 10,000 | 136 | 65.6 | 26.8 | -- | 22.0 | ND | 1.9 | 8.8 | 3.1 | 4.0 | <3.0 | 0.82 J |
| Total Petroleum Hydrocarbons (TPH) by EPA Method 8015 | | | | | | | | | | | | | |
| TPH | NE | -- | 47,600 | 1,690,000 | -- | 741,000 | 275,000 | -- | -- | -- | -- | -- | -- |

All results are in micrograms per liter (ug/L).

Only detected and selected other analytes shown; all others not detected.

- bold shaded** Exceeds Vermont Groundwater Enforcement Standard (VGES)
- bold italics** Reporting limit for non-detection exceeds VGES.
- "--" Not analyzed.
- ND Not detected; reporting limit not shown.
- <X; U Not detected; reporting limit shown.
- NS* Not sampled, free product present.
- NE No standard established.
- J Estimated concentration below reporting limit or due to quality control limitations.
- B Amount in sample >5x (>10x for acetone) amount in method blank; result valid.
- UB Amount in sample <5x (<10x for acetone) amount in method blank; changed to non-detection.
- "/" Analytical result not available.
- † Groundwater Protection Rule and Strategy, 02/05, and Interim Groundwater Quality Standards, Guidance, 08/07.



TABLE 2
Groundwater Samples - Summary of Analytical Results
 Wait Farm Motor Inn
 4805 Main Street, Waitsfield, Vermont

| Sample Identification Sample Date | Vermont Groundwater Enforcement Standard † | MW-3 | | | | | | | | | | |
|--|---|----------|----------|-------------|-------------|-------------|-------------|------------|---------------|----------|-----------------|-----------|
| | | 07/03/02 | 11/21/02 | 02/10/03 | 05/21/03 | 09/25/03 | 03/29/04 | 05/27/05 | 03/29/06 | 05/02/07 | 05/07/08 | 05/18/09 |
| Volatile Organic Compounds (VOCs) by EPA Method 8021B/8260B | | | | | | | | | | | | |
| Acetone | 700 | / | / | / | / | / | / | < 10.0 | < 50.0 | NS* | < 500 | < 50 |
| tertiary-Amyl methyl ether (TAME) | NE | NS* | NS* | -- | -- | -- | -- | < 2.5 | <0.5 J | NS* | <5.0 | <0.50 |
| Benzene | 5.0 | NS* | NS* | 43.6 | 28.0 | 33.2 | 11.3 | 3.2 | <1.0 | NS* | <10.0 | <1.0 |
| n-Butylbenzene | NE | NS* | NS* | -- | -- | -- | -- | 18.9 | 12.4 | NS* | 43.6 | 3.0 |
| sec-Butylbenzene | NE | NS* | NS* | -- | -- | -- | -- | 9.2 | 6.0 | NS* | 20.1 | 1.7 |
| tert-Butylbenzene | NE | / | / | / | / | / | / | < 0.8 | < 1.0 | NS* | < 10 | < 1.0 |
| Chloromethane | 30.0 | / | / | / | / | / | / | < 1.2 | < 2.0 | NS* | < 20.0 | < 2.0 |
| Ethylbenzene | 700.0 | NS* | NS* | 137 | 69.9 | 107 | 80.9 | 47.8 | 29.1 | NS* | 30.9 | 6.1 |
| Isopropylbenzene | NE | NS* | NS* | -- | -- | -- | -- | 15.6 | 11.8 | NS* | 22.9 | 5.5 |
| p-Isopropyltoluene | NE | NS* | NS* | -- | -- | -- | -- | 6.6 | 4.5 | NS* | 12.6 | 1.4 |
| Methyl tert-butyl ether (MTBE) | 40.0 | NS* | NS* | 190 | 46.4 | 42.9 | 32.3 | 19.9 | 5.6 | NS* | <10.0 | <1.0 |
| Naphthalene | 20.0 | NS* | NS* | 494 | 220 | 304 | 226 | 202 | 87.8 J | NS* | 52.2 | 31 |
| n-Propylbenzene | NE | NS* | NS* | -- | -- | -- | -- | 24.6 | 20.0 | NS* | 40.7 | 6.9 |
| 1,1,2,2-Tetrachloroethane | NE | NS* | NS* | -- | -- | -- | -- | < 2.5 | <0.5 | NS* | <5.0 | <0.50 |
| Toluene | 1,000.0 | NS* | NS* | 28.5 | 16.5 | 19.6 | ND | < 3.5 | <1.0 | NS* | <10.0 | <1.0 |
| Trichlorofluoromethane | 2,100.0 | / | / | / | / | / | / | < 0.7 | < 2.0 | NS* | < 20.0 | < 2.0 |
| 1,2,4-Trimethylbenzene | NE | NS* | NS* | 881 | 261 | 323 | 253 | 76.3 | 64.2 | NS* | 105 | 25 |
| 1,3,5-Trimethylbenzene | NE | NS* | NS* | 887 | 96.0 | 135 | 112 | 54.8 | 37.3 | NS* | 70.6 | 15 |
| Total Trimethylbenzenes | 350 | NS* | NS* | 893 | 381 | 507 | 351 | 347.8 | 89.4 | NS* | 223 | 40 |
| Xylenes (total) | 10,000 | NS* | NS* | 940 | 456 | 452 | 388 | 80.4 | 43.1 | NS* | <30.0 | 4.5 |
| Total Petroleum Hydrocarbons (TPH) by EPA Method 8015 | | | | | | | | | | | | |
| TPH | NE | NS* | NS* | 2,830,000 | 16,200 | 184,000 | 59,400 | -- | -- | NS* | -- | -- |

All results are in micrograms per liter (ug/L).

Only detected and selected other analytes shown; all others not detected.

- bold shaded** Exceeds Vermont Groundwater Enforcement Standard (VGES)
- bold italics** Reporting limit for non-detection exceeds VGES.
- "--" Not analyzed.
- ND Not detected; reporting limit not shown.
- <X; U Not detected; reporting limit shown.
- NS* Not sampled, free product present.
- NE No standard established.
- J Estimated concentration below reporting limit or due to quality control limitations.
- B Amount in sample >5x (>10x for acetone) amount in method blank; result valid.
- UB Amount in sample <5x (<10x for acetone) amount in method blank; changed to non-detection.
- "/" Analytical result not available.
- † Groundwater Protection Rule and Strategy, 02/05, and Interim Groundwater Quality Standards, Guidance, 08/07.



TABLE 2
Groundwater Samples - Summary of Analytical Results
 Wait Farm Motor Inn
 4805 Main Street, Waitsfield, Vermont

| Sample Identification Sample Date | Vermont Groundwater Enforcement Standard † | MW-4 | | | | | | | | | | |
|--|---|-------------|-------------|-------------|-------------|-------------|--------------|---------------|----------|----------|-----------|-----------|
| | | 11/21/02 | 02/10/03 | 05/21/03 | 09/25/03 | 03/29/04 | 05/27/05 | 03/29/06 | 05/02/07 | 05/06/08 | 05/18/09 | 06/16/11 |
| Volatile Organic Compounds (VOCs) by EPA Method 8021B/8260B | | | | | | | | | | | | |
| Acetone | 700 | / | / | / | / | / | <200 | <50.0 | <50.0 | <50.0 | <50 | 14 B |
| tertiary-Amyl methyl ether (TAME) | NE | -- | -- | -- | -- | -- | <10.0 | 0.9 J | <0.5 | <0.5 | <0.50 | 2.5 U |
| Benzene | 5.0 | 43.7 | 62.2 | 58.1 | 52.8 | 32.2 | 15.0 | 7.1 | <1.0 | 3.6 | 2.0 | 0.85 J |
| n-Butylbenzene | NE | -- | -- | -- | -- | -- | 28.8 | 15.4 | <1.0 | 17.1 | 13 | 14 |
| sec-Butylbenzene | NE | -- | -- | -- | -- | -- | <12.0 | 9.7 | 1.1 | 12.0 | 8.6 | 9.2 |
| tert-Butylbenzene | NE | / | / | / | / | / | <16.0 | <1.0 | <1.0 | <1.0 | <1.0 | 0.55 J |
| Chloromethane | 30.0 | / | / | / | / | / | <24.0 | <2.0 | <2.0 | <2.0 | <2.0 | 37 |
| Ethylbenzene | 700.0 | 80.8 | 59.0 | 65.9 | 91.4 | 57.8 | 32.6 | 63.2 | <1.0 | 94.8 | 87 | 75 |
| Isopropylbenzene | NE | -- | -- | -- | -- | -- | 35.6 | 22.4 | <1.0 | 28.7 | 23 | 21 |
| p-Isopropyltoluene | NE | -- | -- | -- | -- | -- | 19.8 | 4.8 | <1.0 | 6.2 | 4.7 | 4.8 J |
| Methyl tert-butyl ether (MTBE) | 40.0 | 403 | 273 | 235 | 103 | 86.6 | 41.2 | 23.3 | <1.0 | 6.3 | 3.2 | 2.3 J |
| Naphthalene | 20.0 | 487 | 173 | 204 | 246 | 99.9 | 201 | 104. J | <2.0 | 15.8 | 21 | 5.9 JB |
| n-Propylbenzene | NE | -- | -- | -- | -- | -- | 18.8 | 36.7 | <1.0 | 44.4 | 36 | 33 |
| 1,1,2,2-Tetrachloroethane | NE | -- | -- | -- | -- | -- | <10.0 | <0.5 | <0.5 | <0.5 | <0.50 | 2.5 U |
| Toluene | 1,000.0 | ND | 24.5 | 15.6 | 19.9 | ND | <14.0 | 4.5 | <1.0 | <1.0 | 1.7 | 0.70 J |
| Trichlorofluoromethane | 2,100.0 | / | / | / | / | / | <14.0 | <2.0 | <2.0 | <2.0 | <2.0 | 10 U |
| 1,2,4-Trimethylbenzene | NE | 425 | 180 | 217 | 322 | 99.3 | 30.4 | 107 | <1.0 | 112 | 120 | 120 |
| 1,3,5-Trimethylbenzene | NE | 132 | 82.1 | 81.9 | 124 | 75.0 | 27.8 | 55.8 | <1.0 | 74.7 | 65 | 55 |
| Total Trimethylbenzenes | 350 | 1015 | 592 | 579 | 513 | 276.5 | 392.8 | 187 | <2.0 | 229 | 185 | 175 |
| Xylenes (total) | 10,000 | 269 | 357 | 423 | 536 | 210 | 16.2 | 121.8 | <3.0 | 55.8 | 118 | 70 |
| Total Petroleum Hydrocarbons (TPH) by EPA Method 8015 | | | | | | | | | | | | |
| TPH | NE | 530,000 | 108,000 | 486,000 | 1,495,000 | 21,500 | -- | -- | -- | -- | -- | -- |

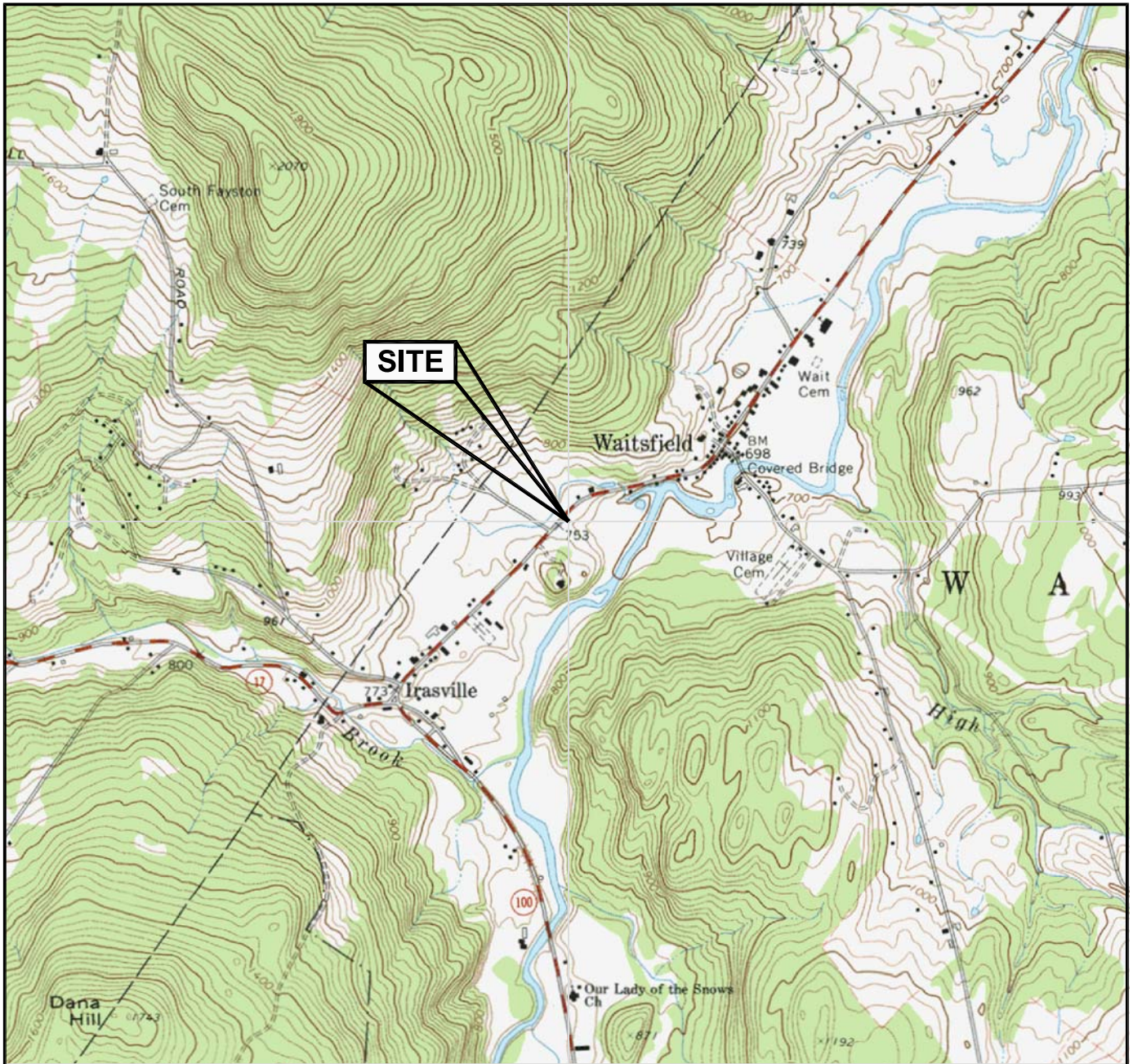
All results are in micrograms per liter (ug/L).

Only detected and selected other analytes shown; all others not detected.

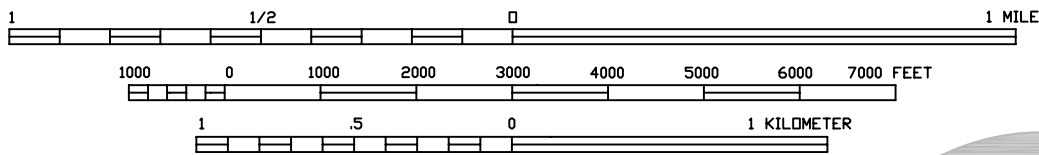
- bold shaded** Exceeds Vermont Groundwater Enforcement Standard (VGES)
- bold italics** Reporting limit for non-detection exceeds VGES.
- "--" Not analyzed.
- ND Not detected; reporting limit not shown.
- <X; U Not detected; reporting limit shown.
- NS* Not sampled, free product present.
- NE No standard established.
- J Estimated concentration below reporting limit or due to quality control limitations.
- B Amount in sample >5x (>10x for acetone) amount in method blank; result valid.
- UB Amount in sample <5x (<10x for acetone) amount in method blank; changed to non-detection.
- "/"/ Analytical result not available.
- † Groundwater Protection Rule and Strategy, 02/05, and Interim Groundwater Quality Standards, Guidance, 08/07.



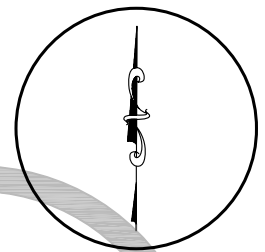
FIGURES



SCALE: 1:24 000



CONTOUR INTERVAL 20 FEET
NATIONAL GEODETIC VERTICAL DATUM OF 1929



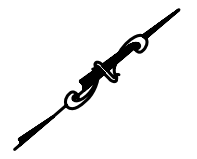
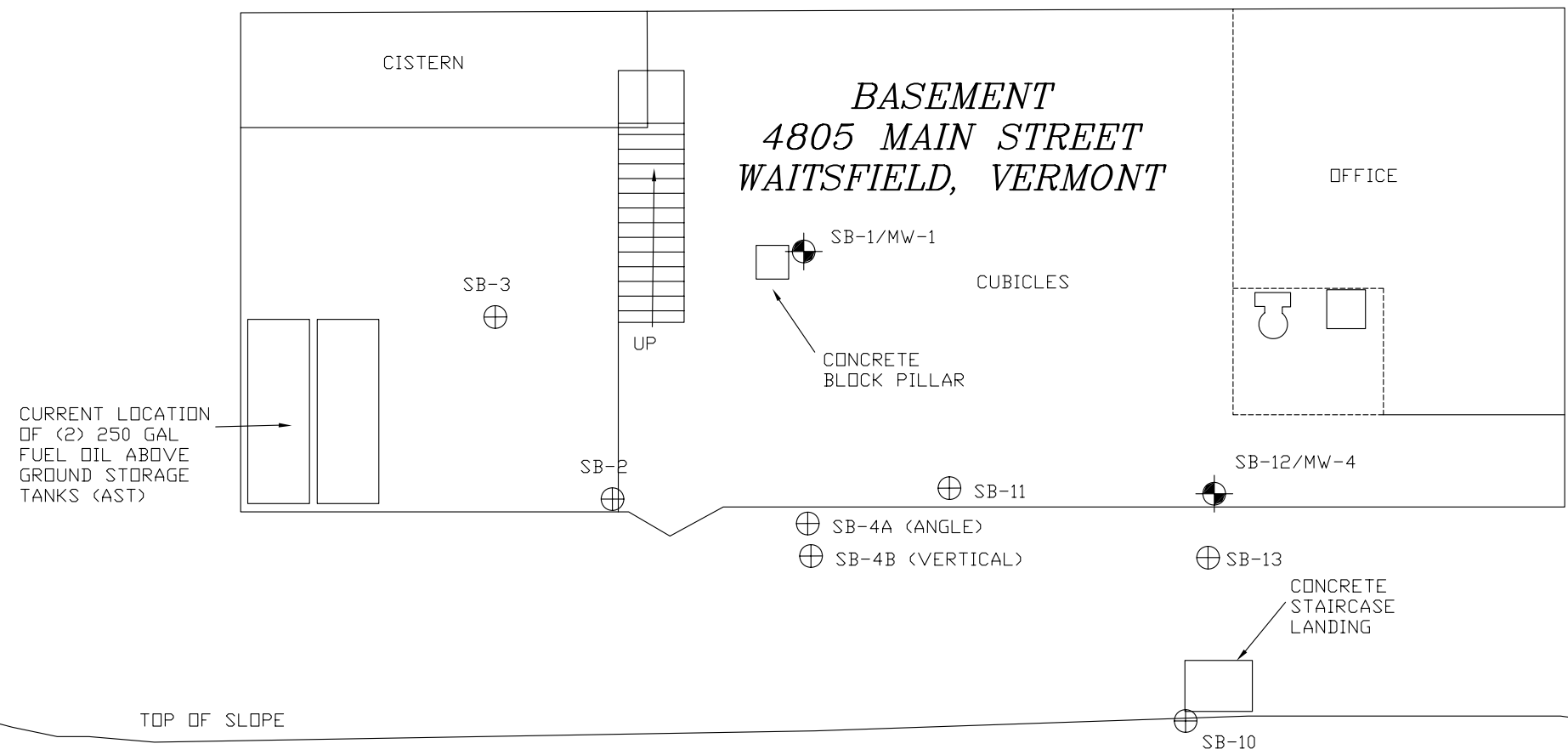
| | | |
|---|---|--------------------------|
| DATE July 8, 2011 | SCALE As shown | FILE Site Loc Map.dwg |
| APPROVED BY RWR | DRAWN BY JDD | REVISED |
| CLIENT Gumbo Group, LLC | JOB NUMBER GUMB0001 | |
| LOCATION Former Wait Farm Motor Inn 4805 Main Street Waitfield, Vermont SMS Site #2002-2982 | MAP SOURCE Waitfield, VT USGS QUAD 1970 | |

Wilcox & Barton INC.
ENVIRONMENTAL AND ENGINEERING SERVICES




SITE LOCATION MAP

Figure 1

**BASEMENT
4805 MAIN STREET
WAITSFIELD, VERMONT**

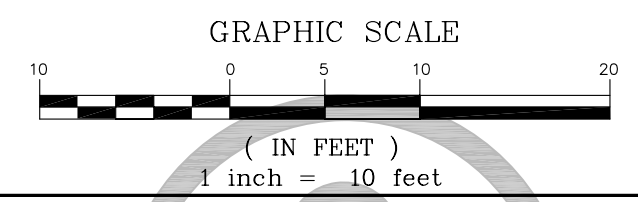


LEGEND

-  SB-12/MW-4 MONITORING WELL
-  SB-9/MW-3 DESTROYED MONITORING WELL
-  SOIL BORING

NOTES

1. ALL LOCATIONS AND DIMENSIONS ARE APPROXIMATE.
2. PLANS BASED ON WILCOX & BARTON, INC. SITE VISIT ON MAY 27, 2005 AND SITE PLANS PRODUCED BY GRIFFIN INTERNATIONAL, INC.
3. THIS PLAN IS NOT A PROFESSIONAL SURVEY AND IS NOT INTENDED FOR SURVEY PURPOSES.



| | | |
|--|-----------------------------------|--------------------------|
| TITLE SITE PLAN | | |
| DATE May 31, 2005 | SCALE 1" = 10' | FILE Waitfarm_Figures |
| APPROVED BY RWR | DRAWN BY JDD | REVISED July 8, 2011 |
| CLIENT Gumbo Group, LLC | JOB NUMBER GUMB0001 | |
| LOCATION Former Wait Farm Motor Inn 4805 Main Street Waitsfield, Vermont SMS Site #2002-2982 | DRAWING NUMBER FIGURE 2 | |

POND

TOP OF SLOPE

APPENDIX A
VT DEC Correspondence



State of Vermont
Department of Environmental Conservation
Waste Management Division
103 South Main Street/West Building
Waterbury, VT 05671-0404
(802) 241-3877
FAX (802) 241-3296
gerold.noyes@state.vt.us

October 9, 2009

PAUL LAVOIE
WAIT FARM MOTOR INN
4805 MAIN STREET
WAITSFIELD, VERMONT 05673

Re: Groundwater monitoring, Wait Farm Motor Inn, SMS Site #2002-2982
Waitsfield, Vermont

Dear Mr. Lavoie:

The Sites Management Section (SMS) has reviewed the October 7, 2009 report prepared by Wilcox & Barton for annual groundwater monitoring conducted on May 18. Naphthalene (31 and 21 $\mu\text{g/L}$) was found above the Vermont Groundwater Enforcement Standards (VGES) in monitor wells MW-3 and -4 respectively. MW-2, -3, and -4 contained other petroleum compounds below the VGES. No contamination was noted in MW-1. Petroleum sheens and odors were noted in the purge water from the monitor wells, although no measurable free product was reported.

Based on these results, the SMS is not requesting additional site investigation or remediation and asks that **biennial** groundwater monitoring be implemented.

We look forward to reviewing the next sample results by **August 15, 2011** and recommendations for future site activity. Please feel free to contact me if I can be of assistance.

Sincerely,

Gerold Noyes, P.E.
Environmental Engineer
Sites Management Section

CC: William Wilcox, Wilcox & Barton (via electronic mail)

Y:\WM_Sites\Sites.Files\2001.2901.to.2002.3000\2002.2982.Wait.Farm.Inn\2009\1009.lavoie.let.doc



APPENDIX B

Site Background Information



SITE BACKGROUND INFORMATION

Wait Farm Motor Inn
4805 Main Street
Waitsfield, Vermont
VT DEC Site #2002-2982

On March 19, 2002, Mr. Paul Lavoie of the Wait Farm Motor Inn retained Griffin International, Inc. (Griffin) to remove a 2,000-gallon fuel oil underground storage tank (UST) from the property. The tank removal was prompted by a leak from a supply line extending to the basement and the presence of petroleum on a pond located near the building.

UST Closure Site Assessment, March 25, 2002, Griffin International, Inc.

On March 25, 2002, Griffin removed the UST and performed a UST closure site assessment. The results of the closure assessment are contained in a letter report to the Vermont Department of Environmental Conservation (VT DEC) dated March 28, 2002. According to the report, the UST and piping exposed during the removal were found to be in good condition with no indication of petroleum leakage or staining on or near the UST. Six confirmatory soil samples from the tank grave were screened for organic vapors with a photoionization detector (PID), revealing a maximum concentration of 13.8 parts per million by volume (ppmv). The cause of the release was attributed to a leaking supply line under the basement. Remedial measures included the recovery of free phase petroleum from the nearby pond.

Initial Report on the Investigation of Subsurface Petroleum Contamination and Product Recovery Efforts, August 28, 2002, Griffin International, Inc.

On August 28, 2002, Griffin produced a report entitled “*Initial Report on the Investigation of Subsurface Petroleum Contamination and Product Recovery Efforts.*” The report summarized site investigative activities implemented for the release, including the installation of 13 soil borings (SB-1 through SB-13) and four one-inch diameter monitoring wells (MW-1 through MW-4) with a hand auger. Soil encountered at the site consisted primarily of silty sand underlain by clay. Soil samples were screened with a PID and selected samples were submitted for laboratory analysis for volatile organic compounds (VOCs) by U.S. Environmental Protection Agency (EPA) Method 8021B. PID measurements revealed petroleum contamination in soil beneath the concrete slab floor of the basement. Soil samples from three borings (SB-2 through SB-4) exhibiting high PID readings were analyzed for VOCs by EPA Method 8021B. The soil analysis revealed petroleum compounds at concentrations above the VT DEC Guidelines for Contaminated Soil and Debris.

Groundwater samples from wells MW-1 and MW-2 were analyzed for VOCs by EPA Method 8021B and found to contain petroleum constituents above the VT DEC Groundwater Enforcement Standards (VGES). A groundwater sample was not collected from monitoring well MW-3 because non-aqueous phase liquid (NAPL) was detected in the well. MW-4 had not yet been installed at the time of the sampling event. Analysis of the on-site potable well for VOCs by EPA Method 524.2 revealed no detection of petroleum constituents.



Griffin recovered product from the pond using sorbent pads and containment booms and from MW-3 using a bailer. It was estimated that approximately 100 gallons of product were recovered. NAPL was measured in MW-3 at a thickness of up to 13.25 inches in July and August of 2002.

Griffin performed a sensitive receptor survey during UST closure and site investigation activities. The pond on the property was noted to overflow to a drainage swale, which ultimately discharges to the Mad River at a location approximately 500 feet to the northeast of the site. Indoor air in the building basement was screened with a PID. The screening revealed a maximum concentration of 2.8 ppmv. The 2.8 ppmv reading was not explained, but it was noted that the basement contained two 275-gallon above-ground storage tanks (ASTs) and an oil-fired boiler.

Griffin recommended groundwater analysis in the fall of 2002 and bi-weekly inspections of the containment area along the pond.

Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination, January 2003, Griffin International, Inc.

Griffin produced a quarterly monitoring report entitled “*Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination*” dated January 2003. The report presented groundwater sampling data from November 21, 2002. Wells MW-1 and MW-3 were found to contain 1/8 inch of NAPL and were, therefore, not sampled. Groundwater samples from MW-2 and MW-4 were analyzed for VOCs by EPA Method 8021B and total petroleum hydrocarbons (TPH) by EPA Method 8015 and found to contain petroleum constituents above VGES. Product was recovered from the monitoring wells that contained NAPL and the absorbents on the pond were monitored and changed as needed.

Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination, March 2003, Griffin International, Inc.

Griffin produced a second quarterly monitoring report entitled “*Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination*” dated March 2003. The report presented groundwater sampling data from February 10, 2003. Groundwater was analyzed for VOCs by EPA Method 8021B and TPH by EPA Method 8015. Petroleum constituents were detected in each well at concentrations above VGES. Product recovery was also performed from monitoring wells containing NAPL and the absorbents on the pond were monitored and changed as needed.

Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination, July 2003, Griffin International, Inc.

Griffin produced a third quarterly monitoring report entitled “*Report on the Quarterly Groundwater Monitoring of Subsurface Petroleum Contamination*” dated July 2003. The report presented groundwater sampling data from May 21, 2003. No NAPL was detected during the sampling event and MW-2 was not sampled because insufficient water was present. Groundwater was analyzed for VOCs by EPA Method 8021B and TPH by EPA Method 8015. Petroleum constituents were detected in each well at concentrations above VGES. Absorbents on the pond and NAPL in the monitoring wells were monitored on a bi-weekly basis during the period. NAPL was detected and recovered from MW-1 and MW-3 on several site visits performed in March and April of 2003. Absorbent booms on the pond were changed as needed. Griffin recommended semi-annual sampling in the spring and fall, monthly NAPL gauging and recovery, and removal of the absorbent materials from the pond. The recommendations were conditionally approved by VT DEC in a letter dated August 1, 2003.

Semi-Annual Groundwater Monitoring Report, October 2003, Griffin International, Inc.

Griffin produced a “*Semi-Annual Groundwater Monitoring Report*” dated October 2003. The report presented groundwater sampling data from September 25, 2003. No NAPL was detected during the sampling event. Groundwater was analyzed for VOCs by EPA Method 8021B and TPH by EPA Method 8015. Petroleum constituents were detected in each well at concentrations above VGES. The monitoring wells were not found to contain NAPL during gauging events in July and September of 2003. The absorbent materials in the pond were removed on September 30, 2003 and removed from the site for disposal. Griffin recommended continuation of the groundwater analysis program on a semi-annual basis, discontinuing TPH analysis, and discontinuing the NAPL gauging and recovery events. VT DEC approved of the Griffin recommendations in a letter dated November 6, 2003.

Semi-Annual Groundwater Monitoring Report, May 2004, Griffin International, Inc.

Griffin produced a second “*Semi-Annual Groundwater Monitoring Report*” dated May 2004. The report presented groundwater sampling data from March 29, 2004. No NAPL was detected during the sampling event. Groundwater was analyzed for VOCs by EPA Method 8021B and TPH by EPA Method 8015. Groundwater was found to contain petroleum constituents in each well at concentrations above VGES. Griffin recommended continuation of the groundwater analysis program on a semi-annual basis, with groundwater analysis for VOCs only.

VT DEC Correspondence, May 21, 2004.

VT DEC produced a letter dated May 21, 2004 providing comments on the Griffin Semi-Annual Groundwater Monitoring Report of May 2004. VT DEC indicated that annual groundwater sampling in the spring was a sufficient scope of work for the project, and requested that TPH sampling be discontinued.

In March 2005, Wilcox & Barton, Inc. was retained to provide environmental consulting services for the site.

Annual Groundwater Monitoring Report, 2005, July 18, 2005, Wilcox & Barton, Inc.

On July 18, 2005, Wilcox & Barton, Inc. produced a report entitled “*Annual Groundwater Monitoring Report, 2005.*” According to the report, an annual groundwater monitoring event was performed on May 27, 2005. Groundwater and product level gauging revealed the presence of 0.02 foot of NAPL in monitoring well MW-1. Prior to this event, NAPL had not been observed on the site since April 2003. Laboratory analysis of groundwater samples from monitoring wells MW-2, MW-3, and MW-4 for VOCs revealed the presence of petroleum constituents in groundwater at concentrations above VGES. Aside from the recurrence of NAPL in MW-1, dissolved contaminant concentrations within each well showed a decreasing trend when compared with historic monitoring data.

The presence of NAPL at the site, as indicated by the 2005 monitoring data, was considered to represent a continuing source of dissolved and vapor phase soil gas contamination. Such conditions were considered to pose a significant threat of impact to surface water and indoor air at the site.

Wilcox & Barton, Inc. recommended increasing the frequency of groundwater monitoring to assess the presence of NAPL in the formation. Groundwater gauging on a quarterly basis during the four seasons of one year was recommended, along with removal of NAPL from the well(s) to ensure the well would continue to accurately reflect dynamic aquifer conditions. Groundwater monitoring for dissolved VOC constituents and reporting was recommended to continue on an annual basis.

Due to the presence of NAPL and potential threats to sensitive receptors, Wilcox & Barton, Inc. recommended that a Corrective Action Feasibility Investigation be performed. It was noted that additional site characterization information may be required to support the completion of the investigation. The required characterization was noted to potentially include the following elements, among others:

- installation of additional monitoring wells to determine the extent of contamination;
- collection and analysis of groundwater samples to determine contaminant distribution and geochemical and groundwater chemistry characteristics indicative of natural attenuation;
- collection of soil vapor samples to assess the potential for migration to indoor air;
- collection and analysis of surface water and/or pond sediment samples;
- surveying of monitoring wells to aid in assessment of contaminant migration rates; and,
- collection and analysis of indoor air samples.

VT DEC Correspondence, February 7, 2006.

On February 7, 2005, VT DEC issued a letter regarding the July 18, 2005, “*Annual Groundwater Monitoring Report, 2005*” prepared by Wilcox & Barton, Inc. VT DEC concurred with the recommendations of the report to increase sampling frequency and product removal, with further

work to be determined based on observed conditions. VT DEC asked that the quarterly monitoring not be initiated until the results of the spring 2006 event were received and reviewed.

Annual Groundwater Monitoring Report, 2006, May 12, 2006, Wilcox & Barton, Inc.

On May 12, 2006, Wilcox & Barton, Inc. produced a report entitled “*Annual Groundwater Monitoring Report, 2006.*” According to the report, an annual groundwater monitoring event was performed on March 29, 2006. Groundwater and product level gauging revealed the presence of 0.02 foot of NAPL in monitoring well MW-1. Laboratory analysis of groundwater samples from monitoring wells MW-2, MW-3, and MW-4 for VOCs revealed the presence of petroleum constituents in groundwater at concentrations above VGES. Aside from the recurrence of NAPL in MW-1, dissolved contaminant concentrations within each well showed a decreasing trend when compared with historic monitoring data.

Wilcox & Barton, Inc. recommended increasing the frequency of groundwater monitoring to assess the presence of NAPL in the formation. Groundwater gauging on a quarterly basis during the four seasons of one year was recommended, along with removal of NAPL from the well(s) to ensure the well would continue to accurately reflect dynamic aquifer conditions. Groundwater monitoring for dissolved VOC constituents and reporting was recommended to continue on an annual basis.

- Due to the presence of NAPL and potential threats to sensitive receptors, Wilcox & Barton, Inc. recommended that a Corrective Action Feasibility Investigation be performed. It was noted that additional site characterization information may be required to support the completion of the investigation.

VT DEC Correspondence, June 1, 2006.

On June 1, 2006, VT DEC issued a letter regarding the May 12, 2006 “*Annual Groundwater Monitoring Report, 2006*” prepared by Wilcox & Barton, Inc. VT DEC did not find the need for a Corrective Action Feasibility Investigation, but did concur with the recommendations to continue annual sampling and quarterly free product removal, with further work to be determined based on observed conditions.

Annual Groundwater Monitoring Report, 2007, June 12, 2007, Wilcox & Barton, Inc.

On June 12, 2007, Wilcox & Barton, Inc. produced a report entitled “*Annual Groundwater Monitoring Report, 2007.*” According to the report, quarterly groundwater and NAPL gauging events were performed on July 21, 2006, October 21, 2006, January 24, 2007, and in conjunction with an annual groundwater monitoring event on May 2, 2007.

The gauging identified the presence of NAPL in monitoring wells MW-1 and MW-3 during the past year of quarterly gauging. The groundwater monitoring event identified the presence of VOCs consistent with fuel oil at concentrations above VGES in samples from wells MW-1 and MW-2.

Wilcox & Barton, Inc. concluded that the current monitoring data indicated that a significant quantity of free-phase fuel oil is present in the subsurface at the site, and represents a continuing source of dissolved and vapor phase soil gas contamination. Such conditions pose a significant threat of impact to surface water and indoor air at the site.

Wilcox & Barton, Inc. recommended the continuation of the quarterly gauging and NAPL recovery for an additional one-year period. Due to the presence of NAPL and potential threats to sensitive receptors, Wilcox & Barton, Inc. continued to recommend that a Corrective Action Feasibility Investigation be performed. Additional site characterization information may be required to support the completion of the investigation and may include the following elements:

- installation of additional monitoring wells to determine the extent of contamination;
- collection and analysis of groundwater samples to determine contaminant distribution and geochemical and groundwater chemistry characteristics indicative of natural attenuation;
- collection of soil vapor samples to assess the potential for migration to indoor air;
- collection and analysis of surface water and/or pond sediment samples;
- surveying of monitoring wells to aid in assessment of contaminant migration rates; and, collection and analysis of indoor air samples.

VT DEC Correspondence, August 1, 2007.

On August 1, 2007, VT DEC issued a letter regarding the *2007 Annual Groundwater Monitoring Report*. The letter referenced the reported free-product recovery results and recommended the discontinuation of free-product recovery.

Annual Groundwater Monitoring Report, 2008, July 14, 2008, Wilcox & Barton, Inc.

On July 14, 2008, Wilcox & Barton submitted a report entitled “Annual Groundwater Monitoring Report, 2008” detailing the monitoring activities and sampling results from 2008. It also included a recommendation for completion of a Corrective Action Feasibility Study and collection of site characterization and exposure pathway data to support that effort. To date, these recommendations have not been accepted by VT DEC.

Annual Groundwater Monitoring Report, 2009, October 7, 2009, Wilcox & Barton, Inc.

On October 7, 2009, Wilcox & Barton submitted a report entitled “Annual Groundwater Monitoring Report, 2009” detailing the monitoring activities and sampling results from 2009. The report again included a recommendation for completion of a Corrective Action Feasibility Study and collection of site characterization and exposure pathway data to support that effort. The VTDEC correspondence dated October 9, 2009, stated further site investigation or remediation would not be required, and biennial sampling and reporting should be implemented.

APPENDIX C

Wilcox & Barton, Inc. Field Sampling Protocol



WILCOX & BARTON, INC.

Protocol for Sampling Groundwater for Volatile Organic Compounds (VOCs) Volatile Petroleum Hydrocarbons (VPH) and Extractable Petroleum Hydrocarbons (EPH)

Groundwater samples may be collected using conventional (standard) techniques or via low-flow techniques depending upon project objectives and data needs. Users are responsible for selecting the appropriate technique and adhering to the protocols outlined below.

STANDARD PURGING AND SAMPLE COLLECTION:

1. Measure depth to static water level and depth to bottom of well.
2. Calculate standing water volume in the well.
3. Using a new or decontaminated bailer, purge either: (1) five well volumes of water from the well, (2) until the well is dry, or (3) until pH, specific conductance and temperature readings stabilize.
4. Transfer the sample directly from the bailer into the appropriate sample container(s).
 - a) VPH and VOC samples – three pre-cleaned, pre-preserved 40-milliliter VOA vials with Teflon septa caps.
 - b) EPH samples - two pre-cleaned 1-liter amber bottles equipped with Teflon-lined screw caps.
5. Pre-preserved sample containers provided by the laboratory are to be used whenever possible. Care must be taken to ensure that the preservative (hydrochloric acid) is not spilled during filling of containers. Sample pH of <2 must be maintained, and can be accomplished by adding 3 to 4 drops of HCl to a 40-ml VOA vial when needed.
6. After filling vials for VOC or VPH samples, invert each vial and tap to liberate potential air bubbles. Inspect to ensure no air bubbles are in the vial.
7. Place samples on ice immediately. Keep refrigerated until delivery to laboratory under chain-of-custody. Samples must be maintained at a temperature of less than 6°C but not frozen.

LOW FLOW PURGING AND SAMPLE COLLECTION (Less than 25 ft to water):

1. Measure depth to static water level and depth to bottom of well.
2. Determine desired sample intake depth based on well construction log and carefully install decontaminated or new/unused sampling tubing and foot valve assembly, minimizing disturbance of the water column:
 - a) Set the intake near the top of the well screen for surficial groundwater monitoring wells; or
 - b) Set the intake at the center of the screened interval for vertical delineation

groundwater monitoring wells.

3. Attach the sampling tubing to clean, unused peristaltic pump tubing using appropriately sized vacuum fittings (male-male hose barb connection or similar fitting).
4. Attach the discharge of the pump to a calibrated flow-through meter capable of measuring, at a minimum, pH, temperature, and specific conductance, and at least one additional parameter such as oxidation-reduction potential, dissolved oxygen, and/or turbidity.
5. Initiate pumping at the lowest possible flow rate and monitor water level drawdown in the well while gradually increasing flow rate. Set the flow rate at the point just before water level drawdown occurs (generally less than 0.5 liter per minute or 0.12 gallons per minute).
6. Observe and record field parameters and water level drawdown initially at 5 to 15 minute intervals. Once stabilization appears to be approaching, increase the observation frequency to every 3 to 5 minutes. Purging is complete when three successive measurements meet the following minimum criteria:
 - pH \pm 0.1
 - Specific Conductivity \pm 0.3%
 - Dissolved Oxygen \pm 10%

If additional parameters are considered, recommended stabilized tolerances are:

- Reduction-Oxidation Potential \pm 10 mV
 - Turbidity \pm 10%
7. Disconnect the flow-through field parameter meter and collect the sample from the pump discharge tubing.
 8. Fill the sample containers directly from the pump discharge tubing, following the same procedures outline in steps 4 through 7 of Standard Purging and Sample Collection above.

DECONTAMINATION:

Use of disposable sampling equipment is preferred. However, if any item of equipment will be introduced into more than one well in a sampling round, the wells should be gauged and/or sampled in order of increasing contamination (*i.e.*, wells anticipated to be less contaminated will be sampled first). After use in each well, the equipment will be decontaminated using, at a minimum, the following sequence:

1. Three rinses with clean water
2. Wash with Alconox (laboratory detergent) and clean water solution
3. Three rinses with clean water
4. Two rinses with distilled water

Similar decontamination procedures should be followed for all equipment introduced into the well, including water level meters, interface probes, and intake tubing. Bailers exposed to free product will not be re-used to sample for dissolved volatile constituents.

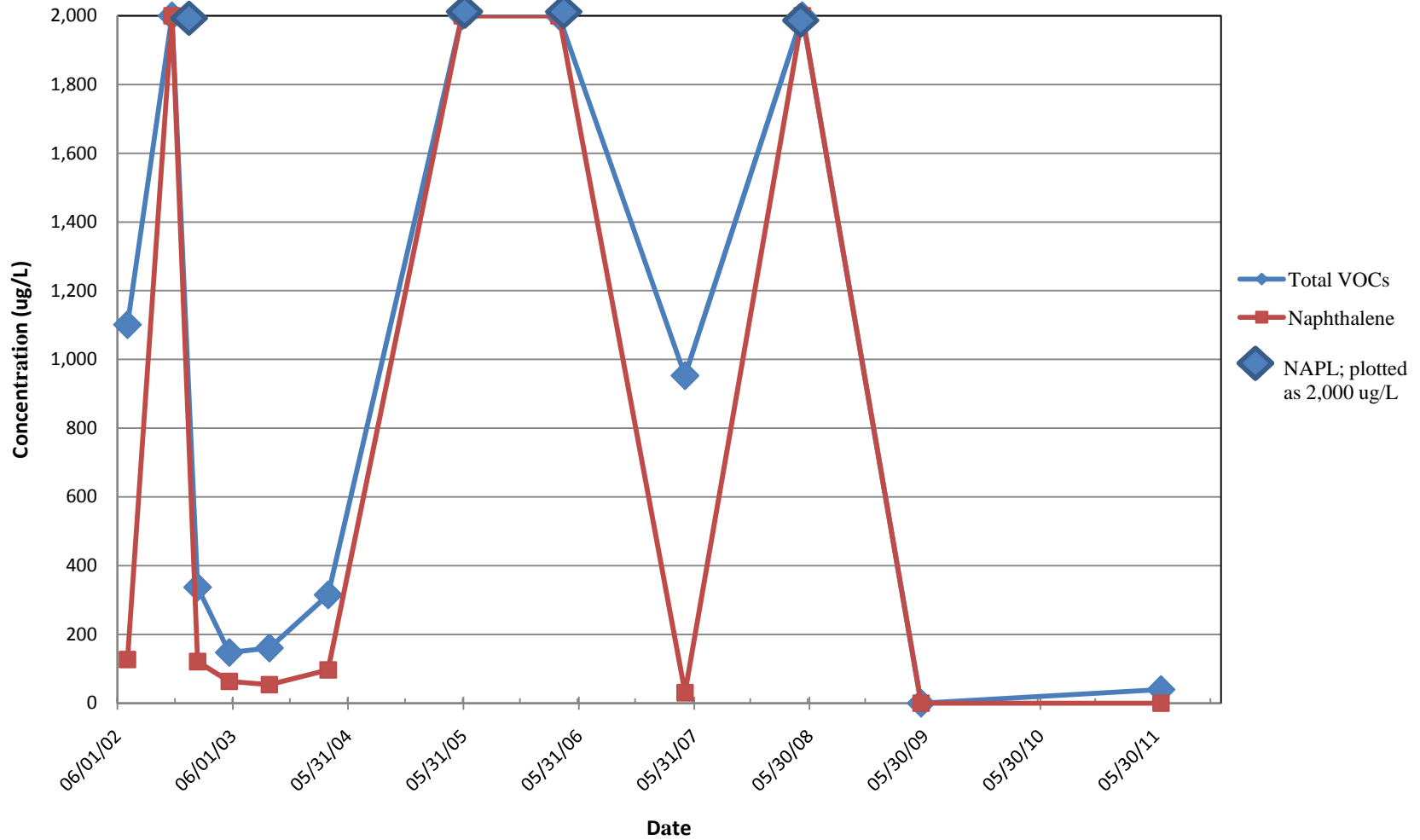
APPENDIX D

Contaminant Trend Graphs



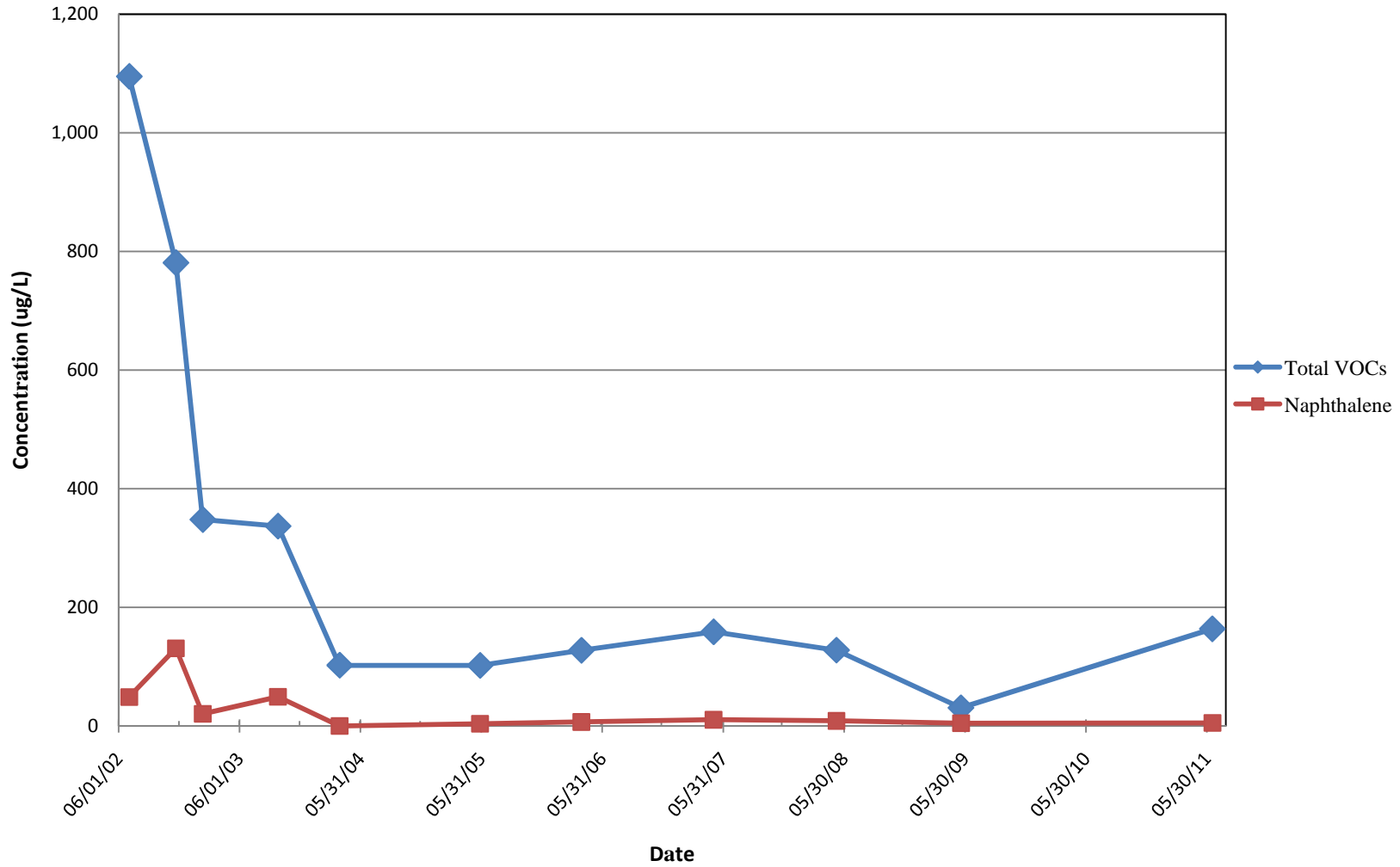
Concentration Trends in MW-1

Former Wait Farm Motor Inn
4805 Main Street, Waitsfield, VT

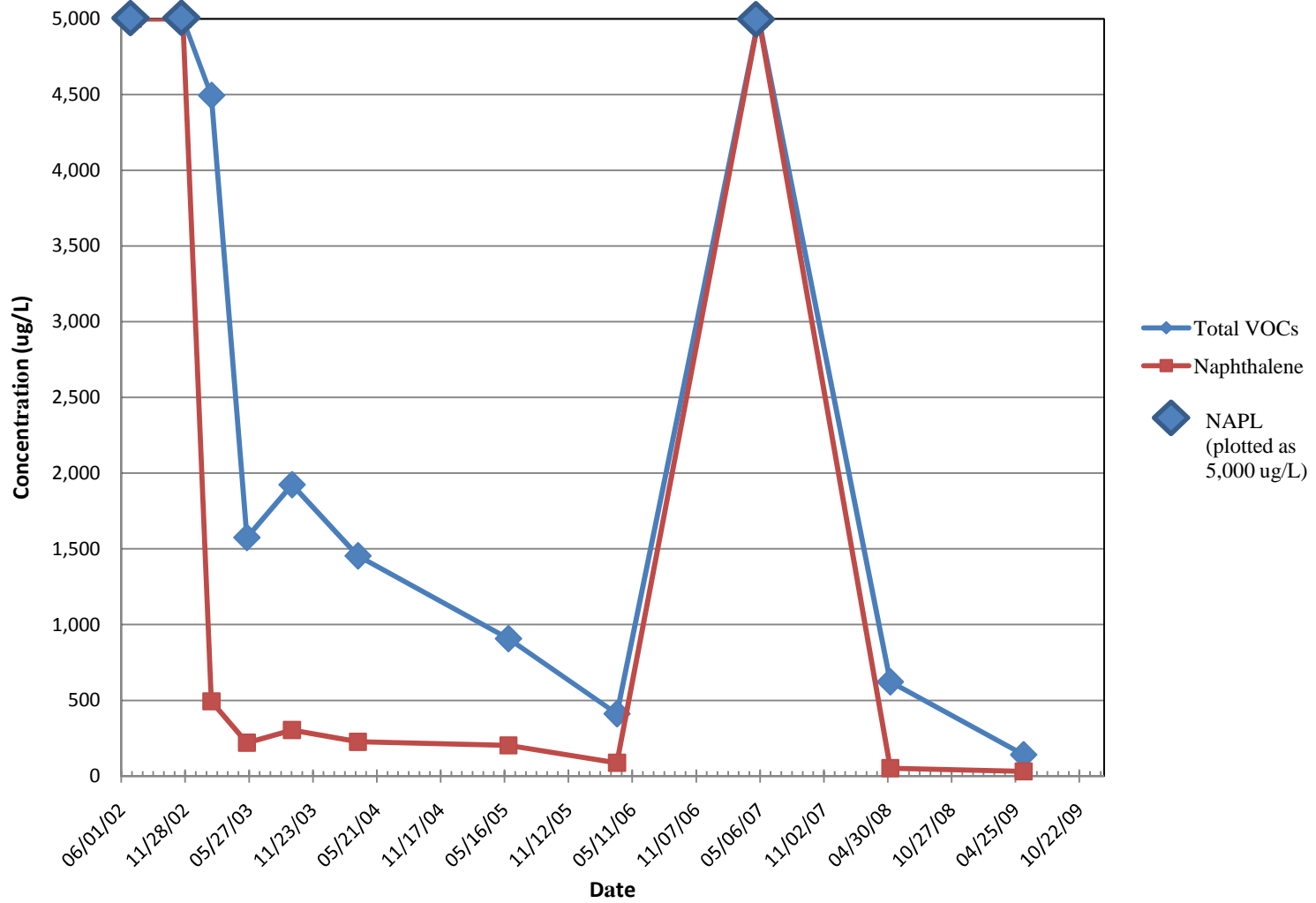


Concentration Trends in MW-2

Former Wait Farm Motor Inn
4805 Main Street, Waitsfield, VT

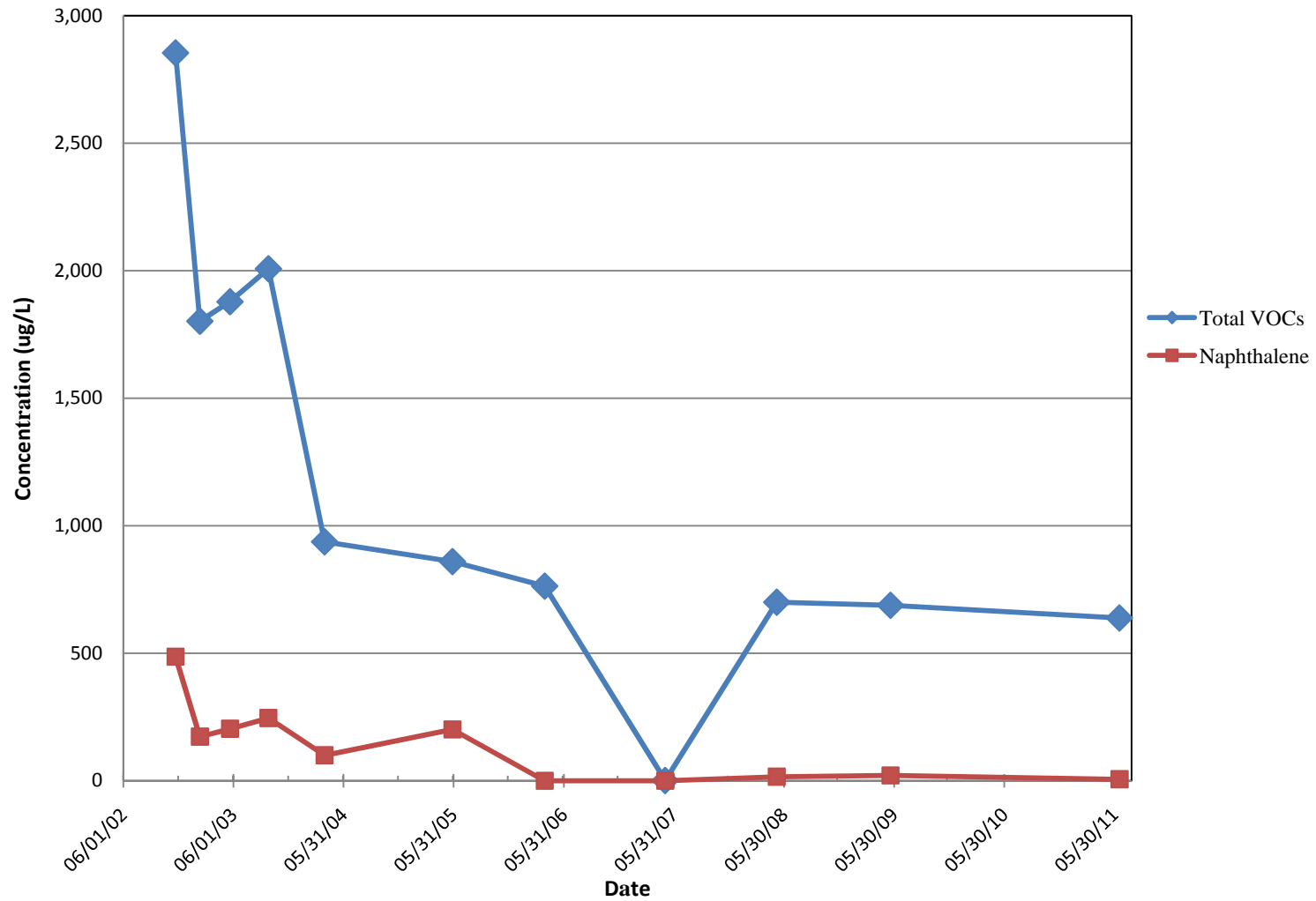


Concentration Trends in MW-3
Former Wait Farm Motor Inn
4805 Main Street, Waitsfield, VT



Concentration Trends in MW-4

Former Wait Farm Motor Inn
4805 Main Street, Waitsfield, VT



APPENDIX E
Laboratory Report



June 28, 2011

James Donaldson
Wilcox & Barton
1115 Route 100B, Suite 200
Moretown, VT 05660

Project Location: 4805 Main St., Waitsfield, VT
Client Job Number:
Project Number: [none]
Laboratory Work Order Number: 11F0756

Enclosed are results of analyses for samples received by the laboratory on June 21, 2011. If you have any questions concerning this report, please feel free to contact me.

Sincerely,



James M. Georgantas
Project Manager

Wilcox & Barton
1115 Route 100B, Suite 200
Moretown, VT 05660
ATTN: James Donaldson

REPORT DATE: 6/28/2011

PURCHASE ORDER NUMBER:

PROJECT NUMBER: [none]

ANALYTICAL SUMMARY

WORK ORDER NUMBER: 11F0756

The results of analyses performed on the following samples submitted to the CON-TEST Analytical Laboratory are found in this report.

PROJECT LOCATION: 4805 Main St., Waitsfield, VT

| FIELD SAMPLE # | LAB ID: | MATRIX | SAMPLE DESCRIPTION | TEST | SUB LAB |
|----------------|------------|--------------|--------------------|--------------|---------|
| MW-1 | 11F0756-01 | Ground Water | | SW-846 8260C | |
| MW-2 | 11F0756-02 | Ground Water | | SW-846 8260C | |
| MW-4 | 11F0756-03 | Ground Water | | SW-846 8260C | |

CASE NARRATIVE SUMMARY

All reported results are within defined laboratory quality control objectives unless listed below or otherwise qualified in this report.

SW-846 8260C

Qualifications:

Laboratory fortified blank/laboratory control sample recovery and duplicate recoveries outside of control limits. Data validation is not affected since all results are "not detected" for associated samples in this batch and bias is on the high side.

Analyte & Samples(s) Qualified:

1,4-Dioxane

B032567-BS1, B032567-BSD1

Either laboratory fortified blank/laboratory control sample or duplicate recovery is outside of control limits, but the other is within limits. RPD between the two LFB/LCS results is within method specified criteria.

Analyte & Samples(s) Qualified:

Dichlorodifluoromethane (Freon 12), Isopropylbenzene (Cumene)

B032642-BSD1, B032567-BSD1

Continuing calibration did not meet method specifications and was biased on the low side for this compound. Increased uncertainty is associated with the reported value which is likely to be biased on the low side.

Analyte & Samples(s) Qualified:

1,2,3-Trichlorobenzene, 1,2,4-Trichlorobenzene, 2,2-Dichloropropane, Naphthalene

11F0756-01[MW-1], 11F0756-02[MW-2], 11F0756-03[MW-4], B032567-BLK1, B032567-BS1, B032567-BSD1, B032642-BLK1, B032642-BS1, B032642-BSD1

Response factor is less than method specified minimum acceptable value. Reduced precision and accuracy are associated with reported result.

Analyte & Samples(s) Qualified:

1,4-Dioxane, tert-Butyl Alcohol (TBA)

11F0756-01[MW-1], 11F0756-02[MW-2], 11F0756-03[MW-4], B032567-BLK1, B032567-BS1, B032567-BSD1, B032642-BLK1, B032642-BS1, B032642-BSD1

Continuing calibration did not meet method specifications and was biased on the high side. Data validation is not affected since sample result was "not detected" for this compound.

Analyte & Samples(s) Qualified:

1,4-Dioxane

B032567-BS1, B032567-BSD1, B032642-BS1, B032642-BSD1

The results of analyses reported only relate to samples submitted to the Con-Test Analytical Laboratory for testing.

I certify that the analyses listed above, unless specifically listed as subcontracted, if any, were performed under my direction according to the approved methodologies listed in this document, and that based upon my inquiry of those individuals immediately responsible for obtaining the information, the material contained in this report is, to the best of my knowledge and belief, accurate and complete.

A handwritten signature in black ink, appearing to read "Daren J. Damboragian", is written over a light gray rectangular background.

Daren J. Damboragian
Laboratory Manager

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-1

Sampled: 6/16/2011 10:00

Sample ID: 11F0756-01

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|------------------------------------|---------|------|-------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Acetone | 2.8 | 50 | 0.54 | µg/L | 1 | J | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Acrylonitrile | ND | 5.0 | 0.51 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| tert-Amyl Methyl Ether (TAME) | ND | 0.50 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Benzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Bromobenzene | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Bromochloromethane | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Bromodichloromethane | ND | 0.50 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Bromoform | ND | 1.0 | 0.25 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Bromomethane | ND | 2.0 | 0.38 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 2-Butanone (MEK) | ND | 20 | 0.41 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| tert-Butyl Alcohol (TBA) | ND | 20 | 3.5 | µg/L | 1 | V-16, U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| n-Butylbenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| sec-Butylbenzene | 0.14 | 1.0 | 0.050 | µg/L | 1 | J | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| tert-Butylbenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| tert-Butyl Ethyl Ether (TBEE) | ND | 0.50 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Carbon Disulfide | ND | 2.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Carbon Tetrachloride | ND | 5.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Chlorobenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Chlorodibromomethane | ND | 1.0 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Chloroethane | ND | 2.0 | 0.33 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Chloroform | ND | 2.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Chloromethane | 38 | 2.0 | 0.13 | µg/L | 1 | | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 2-Chlorotoluene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 4-Chlorotoluene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2-Dibromo-3-chloropropane (DBCP) | ND | 5.0 | 0.48 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2-Dibromoethane (EDB) | ND | 0.50 | 0.14 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Dibromomethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| trans-1,4-Dichloro-2-butene | ND | 2.0 | 0.77 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Dichlorodifluoromethane (Freon 12) | ND | 2.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1-Dichloroethane | ND | 1.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2-Dichloroethane | ND | 1.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1-Dichloroethylene | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| cis-1,2-Dichloroethylene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| trans-1,2-Dichloroethylene | ND | 1.0 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2-Dichloropropane | ND | 1.0 | 0.20 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,3-Dichloropropane | ND | 0.50 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 2,2-Dichloropropane | ND | 1.0 | 0.13 | µg/L | 1 | V-05, U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1-Dichloropropene | ND | 2.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| cis-1,3-Dichloropropene | ND | 0.50 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| trans-1,3-Dichloropropene | ND | 0.50 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-1

Sampled: 6/16/2011 10:00

Sample ID: 11F0756-01

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|------|-------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Diethyl Ether | ND | 2.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Diisopropyl Ether (DIPE) | ND | 0.50 | 0.030 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,4-Dioxane | ND | 50 | 3.5 | µg/L | 1 | V-16, U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Ethylbenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Hexachlorobutadiene | ND | 1.0 | 0.26 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 2-Hexanone (MBK) | ND | 10 | 0.66 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Isopropylbenzene (Cumene) | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| p-Isopropyltoluene (p-Cymene) | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Methyl tert-Butyl Ether (MTBE) | 0.29 | 1.0 | 0.050 | µg/L | 1 | J | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Methylene Chloride | ND | 5.0 | 2.3 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 4-Methyl-2-pentanone (MIBK) | ND | 10 | 0.22 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Naphthalene | ND | 2.0 | 0.21 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| n-Propylbenzene | ND | 1.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Styrene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1,2,2-Tetrachloroethane | ND | 0.50 | 0.18 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Tetrachloroethylene | ND | 1.0 | 0.14 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Tetrahydrofuran | ND | 10 | 1.0 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Toluene | ND | 1.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2,3-Trichlorobenzene | ND | 5.0 | 0.22 | µg/L | 1 | V-05, U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,3,5-Trichlorobenzene | ND | 1.0 | 0.40 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Trichloroethylene | ND | 1.0 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Trichlorofluoromethane (Freon 11) | 1.3 | 2.0 | 0.070 | µg/L | 1 | J | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2,3-Trichloropropane | ND | 2.0 | 0.21 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 1.0 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,2,4-Trimethylbenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| Vinyl Chloride | ND | 2.0 | 0.16 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| m+p Xylene | ND | 2.0 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |
| o-Xylene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/23/11 | 6/23/11 18:32 | MFF |

| Surrogates | % Recovery | Recovery Limits | Flag |
|-----------------------|------------|-----------------|---------------|
| 1,2-Dichloroethane-d4 | 96.2 | 70-130 | 6/23/11 18:32 |
| Toluene-d8 | 100 | 70-130 | 6/23/11 18:32 |
| 4-Bromofluorobenzene | 100 | 70-130 | 6/23/11 18:32 |

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-2

Sampled: 6/16/2011 11:15

Sample ID: 11F0756-02

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|------------------------------------|---------|------|-------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Acetone | 5.7 | 50 | 0.54 | µg/L | 1 | J | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Acrylonitrile | ND | 5.0 | 0.51 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| tert-Amyl Methyl Ether (TAME) | ND | 0.50 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Benzene | 1.2 | 1.0 | 0.050 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Bromobenzene | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Bromochloromethane | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Bromodichloromethane | ND | 0.50 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Bromoform | ND | 1.0 | 0.25 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Bromomethane | ND | 2.0 | 0.38 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 2-Butanone (MEK) | ND | 20 | 0.41 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| tert-Butyl Alcohol (TBA) | ND | 20 | 3.5 | µg/L | 1 | V-16, U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| n-Butylbenzene | 2.1 | 1.0 | 0.050 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| sec-Butylbenzene | 4.5 | 1.0 | 0.050 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| tert-Butylbenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| tert-Butyl Ethyl Ether (TBEE) | ND | 0.50 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Carbon Disulfide | ND | 2.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Carbon Tetrachloride | ND | 5.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Chlorobenzene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Chlorodibromomethane | ND | 1.0 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Chloroethane | ND | 2.0 | 0.33 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Chloroform | ND | 2.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Chloromethane | 110 | 2.0 | 0.13 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 2-Chlorotoluene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 4-Chlorotoluene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2-Dibromo-3-chloropropane (DBCP) | ND | 5.0 | 0.48 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2-Dibromoethane (EDB) | ND | 0.50 | 0.14 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Dibromomethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2-Dichlorobenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,3-Dichlorobenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,4-Dichlorobenzene | ND | 1.0 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| trans-1,4-Dichloro-2-butene | ND | 2.0 | 0.77 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Dichlorodifluoromethane (Freon 12) | ND | 2.0 | 0.040 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1-Dichloroethane | ND | 1.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2-Dichloroethane | ND | 1.0 | 0.090 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1-Dichloroethylene | ND | 1.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| cis-1,2-Dichloroethylene | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| trans-1,2-Dichloroethylene | ND | 1.0 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2-Dichloropropane | ND | 1.0 | 0.20 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,3-Dichloropropane | ND | 0.50 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 2,2-Dichloropropane | ND | 1.0 | 0.13 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1-Dichloropropene | ND | 2.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| cis-1,3-Dichloropropene | ND | 0.50 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| trans-1,3-Dichloropropene | ND | 0.50 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-2

Sampled: 6/16/2011 11:15

Sample ID: 11F0756-02

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|------|-------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Diethyl Ether | ND | 2.0 | 0.10 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Diisopropyl Ether (DIPE) | ND | 0.50 | 0.030 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,4-Dioxane | ND | 50 | 3.5 | µg/L | 1 | V-16, U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Ethylbenzene | 7.0 | 1.0 | 0.050 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Hexachlorobutadiene | ND | 1.0 | 0.26 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 2-Hexanone (MBK) | ND | 10 | 0.66 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Isopropylbenzene (Cumene) | 9.9 | 1.0 | 0.060 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| p-Isopropyltoluene (p-Cymene) | 0.41 | 1.0 | 0.060 | µg/L | 1 | J | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Methyl tert-Butyl Ether (MTBE) | 1.3 | 1.0 | 0.050 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Methylene Chloride | ND | 5.0 | 2.3 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 4-Methyl-2-pentanone (MIBK) | ND | 10 | 0.22 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Naphthalene | 5.0 | 2.0 | 0.21 | µg/L | 1 | V-05 | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| n-Propylbenzene | 8.9 | 1.0 | 0.040 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Styrene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1,2,2-Tetrachloroethane | ND | 0.50 | 0.18 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Tetrachloroethylene | ND | 1.0 | 0.14 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Tetrahydrofuran | ND | 10 | 1.0 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Toluene | 0.34 | 1.0 | 0.040 | µg/L | 1 | J | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2,3-Trichlorobenzene | ND | 5.0 | 0.22 | µg/L | 1 | V-05, U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2,4-Trichlorobenzene | ND | 1.0 | 0.11 | µg/L | 1 | V-05, U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,3,5-Trichlorobenzene | ND | 1.0 | 0.40 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1,1-Trichloroethane | ND | 1.0 | 0.050 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1,2-Trichloroethane | ND | 1.0 | 0.080 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Trichloroethylene | ND | 1.0 | 0.12 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Trichlorofluoromethane (Freon 11) | ND | 2.0 | 0.070 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2,3-Trichloropropane | ND | 2.0 | 0.21 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 1.0 | 0.11 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,2,4-Trimethylbenzene | 6.1 | 1.0 | 0.060 | µg/L | 1 | | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| 1,3,5-Trimethylbenzene | ND | 1.0 | 0.060 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| Vinyl Chloride | ND | 2.0 | 0.16 | µg/L | 1 | U | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| m+p Xylene | 0.52 | 2.0 | 0.070 | µg/L | 1 | J | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |
| o-Xylene | 0.30 | 1.0 | 0.050 | µg/L | 1 | J | SW-846 8260C | 6/24/11 | 6/24/11 15:00 | LBD |

| Surrogates | % Recovery | Recovery Limits | Flag |
|-----------------------|------------|-----------------|---------------|
| 1,2-Dichloroethane-d4 | 95.6 | 70-130 | 6/24/11 15:00 |
| Toluene-d8 | 97.2 | 70-130 | 6/24/11 15:00 |
| 4-Bromofluorobenzene | 99.8 | 70-130 | 6/24/11 15:00 |

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-4

Sampled: 6/16/2011 10:55

Sample ID: 11F0756-03

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|------------------------------------|---------|-----|------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Acetone | 14 | 250 | 2.7 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Acrylonitrile | ND | 25 | 2.6 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| tert-Amyl Methyl Ether (TAME) | ND | 2.5 | 0.55 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Benzene | 0.85 | 5.0 | 0.25 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Bromobenzene | ND | 5.0 | 0.50 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Bromochloromethane | ND | 5.0 | 0.50 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Bromodichloromethane | ND | 2.5 | 0.40 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Bromoform | ND | 5.0 | 1.2 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Bromomethane | ND | 10 | 1.9 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 2-Butanone (MEK) | ND | 100 | 2.0 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| tert-Butyl Alcohol (TBA) | ND | 100 | 18 | µg/L | 5 | V-16, U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| n-Butylbenzene | 14 | 5.0 | 0.25 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| sec-Butylbenzene | 9.2 | 5.0 | 0.25 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| tert-Butylbenzene | 0.55 | 5.0 | 0.25 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| tert-Butyl Ethyl Ether (TBEE) | ND | 2.5 | 0.35 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Carbon Disulfide | ND | 10 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Carbon Tetrachloride | ND | 25 | 0.45 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Chlorobenzene | ND | 5.0 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Chlorodibromomethane | ND | 5.0 | 0.60 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Chloroethane | ND | 10 | 1.6 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Chloroform | ND | 10 | 0.20 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Chloromethane | 37 | 10 | 0.65 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 2-Chlorotoluene | ND | 5.0 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 4-Chlorotoluene | ND | 5.0 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2-Dibromo-3-chloropropane (DBCP) | ND | 25 | 2.4 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2-Dibromoethane (EDB) | ND | 2.5 | 0.70 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Dibromomethane | ND | 5.0 | 0.40 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2-Dichlorobenzene | ND | 5.0 | 0.30 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,3-Dichlorobenzene | ND | 5.0 | 0.30 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,4-Dichlorobenzene | ND | 5.0 | 0.55 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| trans-1,4-Dichloro-2-butene | ND | 10 | 3.8 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Dichlorodifluoromethane (Freon 12) | ND | 10 | 0.20 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1-Dichloroethane | ND | 5.0 | 0.45 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2-Dichloroethane | ND | 5.0 | 0.45 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1-Dichloroethylene | ND | 5.0 | 0.50 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| cis-1,2-Dichloroethylene | ND | 5.0 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| trans-1,2-Dichloroethylene | ND | 5.0 | 0.35 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2-Dichloropropane | ND | 5.0 | 1.0 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,3-Dichloropropane | ND | 2.5 | 0.40 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 2,2-Dichloropropane | ND | 5.0 | 0.65 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1-Dichloropropene | ND | 10 | 0.50 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| cis-1,3-Dichloropropene | ND | 2.5 | 0.35 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| trans-1,3-Dichloropropene | ND | 2.5 | 0.60 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |

Project Location: 4805 Main St., Waitsfield, VT

Sample Description:

Work Order: 11F0756

Date Received: 6/21/2011

Field Sample #: MW-4

Sampled: 6/16/2011 10:55

Sample ID: 11F0756-03

Sample Matrix: Ground Water

Volatile Organic Compounds by GC/MS

| Analyte | Results | RL | DL | Units | Dilution | Flag | Method | Date Prepared | Date/Time Analyzed | Analyst |
|---|---------|-----|------|-------|----------|---------|--------------|---------------|--------------------|---------|
| Diethyl Ether | ND | 10 | 0.50 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Diisopropyl Ether (DIPE) | ND | 2.5 | 0.15 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,4-Dioxane | ND | 250 | 18 | µg/L | 5 | V-16, U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Ethylbenzene | 75 | 5.0 | 0.25 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Hexachlorobutadiene | ND | 5.0 | 1.3 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 2-Hexanone (MBK) | ND | 50 | 3.3 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Isopropylbenzene (Cumene) | 21 | 5.0 | 0.30 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| p-Isopropyltoluene (p-Cymene) | 4.8 | 5.0 | 0.30 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Methyl tert-Butyl Ether (MTBE) | 2.3 | 5.0 | 0.25 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Methylene Chloride | ND | 25 | 11 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 4-Methyl-2-pentanone (MIBK) | ND | 50 | 1.1 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Naphthalene | 5.9 | 10 | 1.0 | µg/L | 5 | V-05, J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| n-Propylbenzene | 33 | 5.0 | 0.20 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Styrene | ND | 5.0 | 0.30 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1,1,2-Tetrachloroethane | ND | 5.0 | 0.40 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1,2,2-Tetrachloroethane | ND | 2.5 | 0.90 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Tetrachloroethylene | ND | 5.0 | 0.70 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Tetrahydrofuran | ND | 50 | 5.0 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Toluene | 0.70 | 5.0 | 0.20 | µg/L | 5 | J | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2,3-Trichlorobenzene | ND | 25 | 1.1 | µg/L | 5 | V-05, U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2,4-Trichlorobenzene | ND | 5.0 | 0.55 | µg/L | 5 | V-05, U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,3,5-Trichlorobenzene | ND | 5.0 | 2.0 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1,1-Trichloroethane | ND | 5.0 | 0.25 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1,2-Trichloroethane | ND | 5.0 | 0.40 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Trichloroethylene | ND | 5.0 | 0.60 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Trichlorofluoromethane (Freon 11) | ND | 10 | 0.35 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2,3-Trichloropropane | ND | 10 | 1.0 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 5.0 | 0.55 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,2,4-Trimethylbenzene | 120 | 5.0 | 0.30 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| 1,3,5-Trimethylbenzene | 55 | 5.0 | 0.30 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| Vinyl Chloride | ND | 10 | 0.80 | µg/L | 5 | U | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| m+p Xylene | 41 | 10 | 0.35 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |
| o-Xylene | 29 | 5.0 | 0.25 | µg/L | 5 | | SW-846 8260C | 6/24/11 | 6/24/11 13:56 | LBD |

| Surrogates | % Recovery | Recovery Limits | Flag |
|-----------------------|------------|-----------------|---------------|
| 1,2-Dichloroethane-d4 | 96.5 | 70-130 | 6/24/11 13:56 |
| Toluene-d8 | 97.9 | 70-130 | 6/24/11 13:56 |
| 4-Bromofluorobenzene | 99.1 | 70-130 | 6/24/11 13:56 |

Sample Extraction Data

Prep Method: SW-846 5030B-SW-846 8260C

| Lab Number [Field ID] | Batch | Initial [mL] | Final [mL] | Date |
|------------------------------|--------------|---------------------|-------------------|-------------|
| 11F0756-01 [MW-1] | B032567 | 5 | 5.00 | 06/23/11 |

Prep Method: SW-846 5030B-SW-846 8260C

| Lab Number [Field ID] | Batch | Initial [mL] | Final [mL] | Date |
|------------------------------|--------------|---------------------|-------------------|-------------|
| 11F0756-02 [MW-2] | B032642 | 5 | 5.00 | 06/24/11 |
| 11F0756-03 [MW-4] | B032642 | 1 | 5.00 | 06/24/11 |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032567 - SW-846 5030B

Blank (B032567-BLK1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|------------------------------------|------|------|------|--|--|--|--|--|--|---------|
| Acetone | 0.82 | 50 | µg/L | | | | | | | J |
| Acrylonitrile | ND | 5.0 | µg/L | | | | | | | U |
| tert-Amyl Methyl Ether (TAME) | ND | 0.50 | µg/L | | | | | | | U |
| Benzene | ND | 1.0 | µg/L | | | | | | | U |
| Bromobenzene | ND | 1.0 | µg/L | | | | | | | U |
| Bromochloromethane | ND | 1.0 | µg/L | | | | | | | U |
| Bromodichloromethane | ND | 0.50 | µg/L | | | | | | | U |
| Bromoform | ND | 1.0 | µg/L | | | | | | | U |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | U |
| 2-Butanone (MEK) | ND | 20 | µg/L | | | | | | | U |
| tert-Butyl Alcohol (TBA) | ND | 20 | µg/L | | | | | | | V-16, U |
| n-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| sec-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| tert-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| tert-Butyl Ethyl Ether (TBEE) | ND | 0.50 | µg/L | | | | | | | U |
| Carbon Disulfide | ND | 2.0 | µg/L | | | | | | | U |
| Carbon Tetrachloride | ND | 5.0 | µg/L | | | | | | | U |
| Chlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| Chlorodibromomethane | ND | 0.50 | µg/L | | | | | | | U |
| Chloroethane | ND | 2.0 | µg/L | | | | | | | U |
| Chloroform | ND | 2.0 | µg/L | | | | | | | U |
| Chloromethane | ND | 2.0 | µg/L | | | | | | | U |
| 2-Chlorotoluene | ND | 1.0 | µg/L | | | | | | | U |
| 4-Chlorotoluene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dibromo-3-chloropropane (DBCP) | ND | 5.0 | µg/L | | | | | | | U |
| 1,2-Dibromoethane (EDB) | ND | 0.50 | µg/L | | | | | | | U |
| Dibromomethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,3-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,4-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| trans-1,4-Dichloro-2-butene | ND | 2.0 | µg/L | | | | | | | U |
| Dichlorodifluoromethane (Freon 12) | ND | 2.0 | µg/L | | | | | | | U |
| 1,1-Dichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| cis-1,2-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| trans-1,2-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichloropropane | ND | 1.0 | µg/L | | | | | | | U |
| 1,3-Dichloropropane | ND | 0.50 | µg/L | | | | | | | U |
| 2,2-Dichloropropane | ND | 1.0 | µg/L | | | | | | | V-05, U |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | U |
| cis-1,3-Dichloropropene | ND | 0.50 | µg/L | | | | | | | U |
| trans-1,3-Dichloropropene | ND | 0.50 | µg/L | | | | | | | U |
| Diethyl Ether | ND | 2.0 | µg/L | | | | | | | U |
| Diisopropyl Ether (DIPE) | ND | 0.50 | µg/L | | | | | | | U |
| 1,4-Dioxane | ND | 50 | µg/L | | | | | | | V-16, U |
| Ethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Hexachlorobutadiene | ND | 0.50 | µg/L | | | | | | | U |
| 2-Hexanone (MBK) | ND | 10 | µg/L | | | | | | | U |
| Isopropylbenzene (Cumene) | ND | 1.0 | µg/L | | | | | | | U |
| p-Isopropyltoluene (p-Cymene) | ND | 1.0 | µg/L | | | | | | | U |
| Methyl tert-Butyl Ether (MTBE) | ND | 1.0 | µg/L | | | | | | | U |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032567 - SW-846 5030B

Blank (B032567-BLK1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|---|------|------|------|------|--|------|--------|--|--|---------|
| Methylene Chloride | ND | 5.0 | µg/L | | | | | | | U |
| 4-Methyl-2-pentanone (MIBK) | ND | 10 | µg/L | | | | | | | U |
| Naphthalene | 0.28 | 2.0 | µg/L | | | | | | | J |
| n-Propylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Styrene | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,2,2-Tetrachloroethane | ND | 0.50 | µg/L | | | | | | | U |
| Tetrachloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | U |
| Toluene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2,3-Trichlorobenzene | 0.61 | 5.0 | µg/L | | | | | | | V-05, J |
| 1,2,4-Trichlorobenzene | 0.20 | 1.0 | µg/L | | | | | | | J |
| 1,3,5-Trichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,1-Trichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,2-Trichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| Trichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| Trichlorofluoromethane (Freon 11) | ND | 2.0 | µg/L | | | | | | | U |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L | | | | | | | U |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 1.0 | µg/L | | | | | | | U |
| 1,2,4-Trimethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,3,5-Trimethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Vinyl Chloride | ND | 2.0 | µg/L | | | | | | | U |
| m+p Xylene | ND | 2.0 | µg/L | | | | | | | U |
| o-Xylene | ND | 1.0 | µg/L | | | | | | | U |
| Surrogate: 1,2-Dichloroethane-d4 | 24.2 | | µg/L | 25.0 | | 96.8 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.3 | | µg/L | 25.0 | | 97.1 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 24.4 | | µg/L | 25.0 | | 97.8 | 70-130 | | | |

LCS (B032567-BS1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|-------------------------------|------|------|------|------|--|------|--------|----|------|---|
| Acetone | 119 | 50 | µg/L | 100 | | 119 | 70-160 | 25 | | † |
| Acrylonitrile | 10.0 | 5.0 | µg/L | 10.0 | | 100 | 70-130 | 25 | | |
| tert-Amyl Methyl Ether (TAME) | 10.5 | 0.50 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |
| Benzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |
| Bromobenzene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | 25 | | |
| Bromochloromethane | 11.2 | 1.0 | µg/L | 10.0 | | 112 | 70-130 | 25 | | |
| Bromodichloromethane | 10.8 | 0.50 | µg/L | 10.0 | | 108 | 70-130 | 25 | | |
| Bromoform | 9.04 | 1.0 | µg/L | 10.0 | | 90.4 | 70-130 | 25 | | |
| Bromomethane | 4.79 | 2.0 | µg/L | 10.0 | | 47.9 | 40-160 | 25 | | † |
| 2-Butanone (MEK) | 113 | 20 | µg/L | 100 | | 113 | 40-160 | 25 | | † |
| tert-Butyl Alcohol (TBA) | 110 | 20 | µg/L | 100 | | 110 | 40-160 | 25 | V-16 | † |
| n-Butylbenzene | 9.52 | 1.0 | µg/L | 10.0 | | 95.2 | 70-130 | 25 | | |
| sec-Butylbenzene | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | 25 | | |
| tert-Butylbenzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |
| tert-Butyl Ethyl Ether (TBEE) | 9.24 | 0.50 | µg/L | 10.0 | | 92.4 | 70-130 | 25 | | |
| Carbon Disulfide | 8.79 | 2.0 | µg/L | 10.0 | | 87.9 | 70-130 | 25 | | |
| Carbon Tetrachloride | 8.85 | 5.0 | µg/L | 10.0 | | 88.5 | 70-130 | 25 | | |
| Chlorobenzene | 11.1 | 1.0 | µg/L | 10.0 | | 111 | 70-130 | 25 | | |
| Chlorodibromomethane | 10.7 | 0.50 | µg/L | 10.0 | | 107 | 70-130 | 25 | | |
| Chloroethane | 8.77 | 2.0 | µg/L | 10.0 | | 87.7 | 70-130 | 25 | | |
| Chloroform | 10.3 | 2.0 | µg/L | 10.0 | | 103 | 70-130 | 25 | | |
| Chloromethane | 6.43 | 2.0 | µg/L | 10.0 | | 64.3 | 40-160 | 25 | | † |
| 2-Chlorotoluene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 25 | | |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---|--------|-----------------|-------|-------------|---------------|--------------|-------------|-----|-----------|--------------------|
| Batch B032567 - SW-846 5030B | | | | | | | | | | |
| LCS (B032567-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 06/23/11 | | | | | | | | | | |
| 4-Chlorotoluene | 11.4 | 1.0 | µg/L | 10.0 | | 114 | 70-130 | | 25 | |
| 1,2-Dibromo-3-chloropropane (DBCP) | 9.24 | 5.0 | µg/L | 10.0 | | 92.4 | 70-130 | | 25 | |
| 1,2-Dibromoethane (EDB) | 11.3 | 0.50 | µg/L | 10.0 | | 113 | 70-130 | | 25 | |
| Dibromomethane | 11.4 | 1.0 | µg/L | 10.0 | | 114 | 70-130 | | 25 | |
| 1,2-Dichlorobenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | | 25 | |
| 1,3-Dichlorobenzene | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | | 25 | |
| 1,4-Dichlorobenzene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | | 25 | |
| trans-1,4-Dichloro-2-butene | 8.06 | 2.0 | µg/L | 10.0 | | 80.6 | 70-130 | | 25 | |
| Dichlorodifluoromethane (Freon 12) | 4.90 | 2.0 | µg/L | 10.0 | | 49.0 | 40-160 | | 25 | † |
| 1,1-Dichloroethane | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | | 25 | |
| 1,2-Dichloroethane | 11.1 | 1.0 | µg/L | 10.0 | | 111 | 70-130 | | 25 | |
| 1,1-Dichloroethylene | 9.50 | 1.0 | µg/L | 10.0 | | 95.0 | 70-130 | | 25 | |
| cis-1,2-Dichloroethylene | 9.85 | 1.0 | µg/L | 10.0 | | 98.5 | 70-130 | | 25 | |
| trans-1,2-Dichloroethylene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | | 25 | |
| 1,2-Dichloropropane | 10.8 | 1.0 | µg/L | 10.0 | | 108 | 70-130 | | 25 | |
| 1,3-Dichloropropane | 11.0 | 0.50 | µg/L | 10.0 | | 110 | 70-130 | | 25 | |
| 2,2-Dichloropropane | 6.77 | 1.0 | µg/L | 10.0 | | 67.7 | 40-130 | | 25 | V-05 † |
| 1,1-Dichloropropene | 10.2 | 2.0 | µg/L | 10.0 | | 102 | 70-130 | | 25 | |
| cis-1,3-Dichloropropene | 8.57 | 0.50 | µg/L | 10.0 | | 85.7 | 70-130 | | 25 | |
| trans-1,3-Dichloropropene | 9.15 | 0.50 | µg/L | 10.0 | | 91.5 | 70-130 | | 25 | |
| Diethyl Ether | 10.6 | 2.0 | µg/L | 10.0 | | 106 | 70-130 | | 25 | |
| Diisopropyl Ether (DIPE) | 9.93 | 0.50 | µg/L | 10.0 | | 99.3 | 70-130 | | 25 | |
| 1,4-Dioxane | 155 | 50 | µg/L | 100 | | 155 * | 40-130 | | 50 | L-02, V-16, V-20 † |
| Ethylbenzene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | | 25 | |
| Hexachlorobutadiene | 9.74 | 0.50 | µg/L | 10.0 | | 97.4 | 70-130 | | 25 | |
| 2-Hexanone (MBK) | 117 | 10 | µg/L | 100 | | 117 | 70-160 | | 25 | † |
| Isopropylbenzene (Cumene) | 13.0 | 1.0 | µg/L | 10.0 | | 130 | 70-130 | | 25 | |
| p-Isopropyltoluene (p-Cymene) | 9.92 | 1.0 | µg/L | 10.0 | | 99.2 | 70-130 | | 25 | |
| Methyl tert-Butyl Ether (MTBE) | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Methylene Chloride | 10.5 | 5.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| 4-Methyl-2-pentanone (MIBK) | 116 | 10 | µg/L | 100 | | 116 | 70-160 | | 25 | † |
| Naphthalene | 8.95 | 2.0 | µg/L | 10.0 | | 89.5 | 40-130 | | 25 | † |
| n-Propylbenzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Styrene | 11.6 | 1.0 | µg/L | 10.0 | | 116 | 70-130 | | 25 | |
| 1,1,1,2-Tetrachloroethane | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | | 25 | |
| 1,1,2,2-Tetrachloroethane | 10.1 | 0.50 | µg/L | 10.0 | | 101 | 70-130 | | 25 | |
| Tetrachloroethylene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Tetrahydrofuran | 9.35 | 10 | µg/L | 10.0 | | 93.5 | 70-130 | | 25 | J |
| Toluene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | | 25 | |
| 1,2,3-Trichlorobenzene | 8.11 | 5.0 | µg/L | 10.0 | | 81.1 | 70-130 | | 25 | V-05 |
| 1,2,4-Trichlorobenzene | 8.39 | 1.0 | µg/L | 10.0 | | 83.9 | 70-130 | | 25 | |
| 1,3,5-Trichlorobenzene | 8.91 | 1.0 | µg/L | 10.0 | | 89.1 | 70-130 | | 25 | |
| 1,1,1-Trichloroethane | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | | 25 | |
| 1,1,2-Trichloroethane | 11.1 | 1.0 | µg/L | 10.0 | | 111 | 70-130 | | 25 | |
| Trichloroethylene | 11.2 | 1.0 | µg/L | 10.0 | | 112 | 70-130 | | 25 | |
| Trichlorofluoromethane (Freon 11) | 7.87 | 2.0 | µg/L | 10.0 | | 78.7 | 70-130 | | 25 | |
| 1,2,3-Trichloropropane | 9.69 | 2.0 | µg/L | 10.0 | | 96.9 | 70-130 | | 25 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 8.74 | 1.0 | µg/L | 10.0 | | 87.4 | 70-130 | | 25 | |
| 1,2,4-Trimethylbenzene | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | | 25 | |
| 1,3,5-Trimethylbenzene | 10.8 | 1.0 | µg/L | 10.0 | | 108 | 70-130 | | 25 | |
| Vinyl Chloride | 7.60 | 2.0 | µg/L | 10.0 | | 76.0 | 40-160 | | 25 | † |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032567 - SW-846 5030B

LCS (B032567-BS1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|----------------------------------|------|-----|------|------|--|------|--------|--|----|--|
| m+p Xylene | 23.0 | 2.0 | µg/L | 20.0 | | 115 | 70-130 | | 25 | |
| o-Xylene | 11.8 | 1.0 | µg/L | 10.0 | | 118 | 70-130 | | 25 | |
| Surrogate: 1,2-Dichloroethane-d4 | 23.8 | | µg/L | 25.0 | | 95.2 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.2 | | µg/L | 25.0 | | 97.0 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 25.3 | | µg/L | 25.0 | | 101 | 70-130 | | | |

LCS Dup (B032567-BSD1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|------------------------------------|------|------|------|------|--|------|--------|-------|----|--------|
| Acetone | 146 | 50 | µg/L | 100 | | 146 | 70-160 | 20.1 | 25 | † |
| Acrylonitrile | 10.4 | 5.0 | µg/L | 10.0 | | 104 | 70-130 | 2.94 | 25 | |
| tert-Amyl Methyl Ether (TAME) | 10.7 | 0.50 | µg/L | 10.0 | | 107 | 70-130 | 1.51 | 25 | |
| Benzene | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | 1.42 | 25 | |
| Bromobenzene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 0.641 | 25 | |
| Bromochloromethane | 11.4 | 1.0 | µg/L | 10.0 | | 114 | 70-130 | 1.95 | 25 | |
| Bromodichloromethane | 10.9 | 0.50 | µg/L | 10.0 | | 109 | 70-130 | 0.831 | 25 | |
| Bromoform | 9.06 | 1.0 | µg/L | 10.0 | | 90.6 | 70-130 | 0.221 | 25 | |
| Bromomethane | 5.10 | 2.0 | µg/L | 10.0 | | 51.0 | 40-160 | 6.27 | 25 | † |
| 2-Butanone (MEK) | 121 | 20 | µg/L | 100 | | 121 | 40-160 | 6.30 | 25 | † |
| tert-Butyl Alcohol (TBA) | 107 | 20 | µg/L | 100 | | 107 | 40-160 | 3.13 | 25 | V-16 † |
| n-Butylbenzene | 9.94 | 1.0 | µg/L | 10.0 | | 99.4 | 70-130 | 4.32 | 25 | |
| sec-Butylbenzene | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 3.15 | 25 | |
| tert-Butylbenzene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 5.11 | 25 | |
| tert-Butyl Ethyl Ether (TBEE) | 9.26 | 0.50 | µg/L | 10.0 | | 92.6 | 70-130 | 0.216 | 25 | |
| Carbon Disulfide | 8.38 | 2.0 | µg/L | 10.0 | | 83.8 | 70-130 | 4.78 | 25 | |
| Carbon Tetrachloride | 9.14 | 5.0 | µg/L | 10.0 | | 91.4 | 70-130 | 3.22 | 25 | |
| Chlorobenzene | 11.4 | 1.0 | µg/L | 10.0 | | 114 | 70-130 | 2.85 | 25 | |
| Chlorodibromomethane | 11.0 | 0.50 | µg/L | 10.0 | | 110 | 70-130 | 2.58 | 25 | |
| Chloroethane | 8.49 | 2.0 | µg/L | 10.0 | | 84.9 | 70-130 | 3.24 | 25 | |
| Chloroform | 10.6 | 2.0 | µg/L | 10.0 | | 106 | 70-130 | 3.34 | 25 | |
| Chloromethane | 6.28 | 2.0 | µg/L | 10.0 | | 62.8 | 40-160 | 2.36 | 25 | † |
| 2-Chlorotoluene | 11.2 | 1.0 | µg/L | 10.0 | | 112 | 70-130 | 1.62 | 25 | |
| 4-Chlorotoluene | 11.6 | 1.0 | µg/L | 10.0 | | 116 | 70-130 | 1.83 | 25 | |
| 1,2-Dibromo-3-chloropropane (DBCP) | 9.34 | 5.0 | µg/L | 10.0 | | 93.4 | 70-130 | 1.08 | 25 | |
| 1,2-Dibromoethane (EDB) | 11.3 | 0.50 | µg/L | 10.0 | | 113 | 70-130 | 0.177 | 25 | |
| Dibromomethane | 11.6 | 1.0 | µg/L | 10.0 | | 116 | 70-130 | 1.82 | 25 | |
| 1,2-Dichlorobenzene | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | 2.74 | 25 | |
| 1,3-Dichlorobenzene | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | 0.749 | 25 | |
| 1,4-Dichlorobenzene | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | 1.36 | 25 | |
| trans-1,4-Dichloro-2-butene | 8.31 | 2.0 | µg/L | 10.0 | | 83.1 | 70-130 | 3.05 | 25 | |
| Dichlorodifluoromethane (Freon 12) | 4.79 | 2.0 | µg/L | 10.0 | | 47.9 | 40-160 | 2.27 | 25 | † |
| 1,1-Dichloroethane | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 4.24 | 25 | |
| 1,2-Dichloroethane | 11.6 | 1.0 | µg/L | 10.0 | | 116 | 70-130 | 4.39 | 25 | |
| 1,1-Dichloroethylene | 9.49 | 1.0 | µg/L | 10.0 | | 94.9 | 70-130 | 0.105 | 25 | |
| cis-1,2-Dichloroethylene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 3.00 | 25 | |
| trans-1,2-Dichloroethylene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 0.915 | 25 | |
| 1,2-Dichloropropane | 11.1 | 1.0 | µg/L | 10.0 | | 111 | 70-130 | 3.01 | 25 | |
| 1,3-Dichloropropane | 11.2 | 0.50 | µg/L | 10.0 | | 112 | 70-130 | 2.52 | 25 | |
| 2,2-Dichloropropane | 7.17 | 1.0 | µg/L | 10.0 | | 71.7 | 40-130 | 5.74 | 25 | V-05 † |
| 1,1-Dichloropropene | 10.4 | 2.0 | µg/L | 10.0 | | 104 | 70-130 | 1.36 | 25 | |
| cis-1,3-Dichloropropene | 9.05 | 0.50 | µg/L | 10.0 | | 90.5 | 70-130 | 5.45 | 25 | |
| trans-1,3-Dichloropropene | 9.48 | 0.50 | µg/L | 10.0 | | 94.8 | 70-130 | 3.54 | 25 | |
| Diethyl Ether | 11.1 | 2.0 | µg/L | 10.0 | | 111 | 70-130 | 5.08 | 25 | |
| Diisopropyl Ether (DIPE) | 10.0 | 0.50 | µg/L | 10.0 | | 100 | 70-130 | 0.702 | 25 | |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032567 - SW-846 5030B

LCS Dup (B032567-BSD1)

Prepared & Analyzed: 06/23/11

| | | | | | | | | | | |
|---|------|------|------|------|--|------|----------|-------|----|----------------------|
| 1,4-Dioxane | 141 | 50 | µg/L | 100 | | 141 | * 40-130 | 9.00 | 50 | L-02, V-16, V-20 † ‡ |
| Ethylbenzene | 11.3 | 1.0 | µg/L | 10.0 | | 113 | 70-130 | 2.06 | 25 | |
| Hexachlorobutadiene | 10.1 | 0.50 | µg/L | 10.0 | | 101 | 70-130 | 3.73 | 25 | |
| 2-Hexanone (MBK) | 121 | 10 | µg/L | 100 | | 121 | 70-160 | 2.74 | 25 | † |
| Isopropylbenzene (Cumene) | 13.3 | 1.0 | µg/L | 10.0 | | 133 | * 70-130 | 2.05 | 25 | L-07 |
| p-Isopropyltoluene (p-Cymene) | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 4.34 | 25 | |
| Methyl tert-Butyl Ether (MTBE) | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 0.947 | 25 | |
| Methylene Chloride | 10.6 | 5.0 | µg/L | 10.0 | | 106 | 70-130 | 0.662 | 25 | |
| 4-Methyl-2-pentanone (MIBK) | 117 | 10 | µg/L | 100 | | 117 | 70-160 | 0.797 | 25 | † |
| Naphthalene | 9.10 | 2.0 | µg/L | 10.0 | | 91.0 | 40-130 | 1.66 | 25 | † |
| n-Propylbenzene | 10.8 | 1.0 | µg/L | 10.0 | | 108 | 70-130 | 2.53 | 25 | |
| Styrene | 11.8 | 1.0 | µg/L | 10.0 | | 118 | 70-130 | 2.05 | 25 | |
| 1,1,1,2-Tetrachloroethane | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 0.376 | 25 | |
| 1,1,2,2-Tetrachloroethane | 10.1 | 0.50 | µg/L | 10.0 | | 101 | 70-130 | 0.494 | 25 | |
| Tetrachloroethylene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 4.64 | 25 | |
| Tetrahydrofuran | 8.44 | 10 | µg/L | 10.0 | | 84.4 | 70-130 | 10.2 | 25 | J |
| Toluene | 11.1 | 1.0 | µg/L | 10.0 | | 111 | 70-130 | 2.09 | 25 | |
| 1,2,3-Trichlorobenzene | 8.39 | 5.0 | µg/L | 10.0 | | 83.9 | 70-130 | 3.39 | 25 | V-05 |
| 1,2,4-Trichlorobenzene | 8.45 | 1.0 | µg/L | 10.0 | | 84.5 | 70-130 | 0.713 | 25 | |
| 1,3,5-Trichlorobenzene | 8.93 | 1.0 | µg/L | 10.0 | | 89.3 | 70-130 | 0.224 | 25 | |
| 1,1,1-Trichloroethane | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 1.54 | 25 | |
| 1,1,2-Trichloroethane | 11.2 | 1.0 | µg/L | 10.0 | | 112 | 70-130 | 0.897 | 25 | |
| Trichloroethylene | 11.4 | 1.0 | µg/L | 10.0 | | 114 | 70-130 | 1.59 | 25 | |
| Trichlorofluoromethane (Freon 11) | 8.08 | 2.0 | µg/L | 10.0 | | 80.8 | 70-130 | 2.63 | 25 | |
| 1,2,3-Trichloropropane | 9.83 | 2.0 | µg/L | 10.0 | | 98.3 | 70-130 | 1.43 | 25 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 9.06 | 1.0 | µg/L | 10.0 | | 90.6 | 70-130 | 3.60 | 25 | |
| 1,2,4-Trimethylbenzene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 2.21 | 25 | |
| 1,3,5-Trimethylbenzene | 11.2 | 1.0 | µg/L | 10.0 | | 112 | 70-130 | 3.27 | 25 | |
| Vinyl Chloride | 7.38 | 2.0 | µg/L | 10.0 | | 73.8 | 40-160 | 2.94 | 25 | † |
| m+p Xylene | 23.4 | 2.0 | µg/L | 20.0 | | 117 | 70-130 | 1.81 | 25 | |
| o-Xylene | 11.9 | 1.0 | µg/L | 10.0 | | 119 | 70-130 | 1.10 | 25 | |
| Surrogate: 1,2-Dichloroethane-d4 | 23.7 | | µg/L | 25.0 | | 94.8 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.7 | | µg/L | 25.0 | | 98.7 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 25.2 | | µg/L | 25.0 | | 101 | 70-130 | | | |

Batch B032642 - SW-846 5030B

Blank (B032642-BLK1)

Prepared & Analyzed: 06/24/11

| | | | | | | | | | | |
|-------------------------------|------|------|------|--|--|--|--|--|--|---------|
| Acetone | 0.80 | 50 | µg/L | | | | | | | J |
| Acrylonitrile | ND | 5.0 | µg/L | | | | | | | U |
| tert-Amyl Methyl Ether (TAME) | ND | 0.50 | µg/L | | | | | | | U |
| Benzene | ND | 1.0 | µg/L | | | | | | | U |
| Bromobenzene | ND | 1.0 | µg/L | | | | | | | U |
| Bromochloromethane | ND | 1.0 | µg/L | | | | | | | U |
| Bromodichloromethane | ND | 0.50 | µg/L | | | | | | | U |
| Bromoform | ND | 1.0 | µg/L | | | | | | | U |
| Bromomethane | ND | 2.0 | µg/L | | | | | | | U |
| 2-Butanone (MEK) | ND | 20 | µg/L | | | | | | | U |
| tert-Butyl Alcohol (TBA) | ND | 20 | µg/L | | | | | | | V-16, U |
| n-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| sec-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| tert-Butylbenzene | ND | 1.0 | µg/L | | | | | | | U |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032642 - SW-846 5030B

Blank (B032642-BLK1)

Prepared & Analyzed: 06/24/11

| | | | | | | | | | | |
|------------------------------------|------|------|------|--|--|--|--|--|--|---------|
| tert-Butyl Ethyl Ether (TBEE) | ND | 0.50 | µg/L | | | | | | | U |
| Carbon Disulfide | ND | 2.0 | µg/L | | | | | | | U |
| Carbon Tetrachloride | ND | 5.0 | µg/L | | | | | | | U |
| Chlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| Chlorodibromomethane | ND | 0.50 | µg/L | | | | | | | U |
| Chloroethane | ND | 2.0 | µg/L | | | | | | | U |
| Chloroform | ND | 2.0 | µg/L | | | | | | | U |
| Chloromethane | ND | 2.0 | µg/L | | | | | | | U |
| 2-Chlorotoluene | ND | 1.0 | µg/L | | | | | | | U |
| 4-Chlorotoluene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dibromo-3-chloropropane (DBCP) | ND | 5.0 | µg/L | | | | | | | U |
| 1,2-Dibromoethane (EDB) | ND | 0.50 | µg/L | | | | | | | U |
| Dibromomethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,3-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,4-Dichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| trans-1,4-Dichloro-2-butene | ND | 2.0 | µg/L | | | | | | | U |
| Dichlorodifluoromethane (Freon 12) | ND | 2.0 | µg/L | | | | | | | U |
| 1,1-Dichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| cis-1,2-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| trans-1,2-Dichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2-Dichloropropane | ND | 1.0 | µg/L | | | | | | | U |
| 1,3-Dichloropropane | ND | 0.50 | µg/L | | | | | | | U |
| 2,2-Dichloropropane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1-Dichloropropene | ND | 2.0 | µg/L | | | | | | | U |
| cis-1,3-Dichloropropene | ND | 0.50 | µg/L | | | | | | | U |
| trans-1,3-Dichloropropene | ND | 0.50 | µg/L | | | | | | | U |
| Diethyl Ether | ND | 2.0 | µg/L | | | | | | | U |
| Diisopropyl Ether (DIPE) | ND | 0.50 | µg/L | | | | | | | U |
| 1,4-Dioxane | ND | 50 | µg/L | | | | | | | V-16, U |
| Ethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Hexachlorobutadiene | ND | 0.50 | µg/L | | | | | | | U |
| 2-Hexanone (MBK) | ND | 10 | µg/L | | | | | | | U |
| Isopropylbenzene (Cumene) | ND | 1.0 | µg/L | | | | | | | U |
| p-Isopropyltoluene (p-Cymene) | ND | 1.0 | µg/L | | | | | | | U |
| Methyl tert-Butyl Ether (MTBE) | ND | 1.0 | µg/L | | | | | | | U |
| Methylene Chloride | ND | 5.0 | µg/L | | | | | | | U |
| 4-Methyl-2-pentanone (MIBK) | ND | 10 | µg/L | | | | | | | U |
| Naphthalene | 0.26 | 2.0 | µg/L | | | | | | | V-05, J |
| n-Propylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Styrene | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,1,2-Tetrachloroethane | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,1,2,2-Tetrachloroethane | ND | 0.50 | µg/L | | | | | | | U |
| Tetrachloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| Tetrahydrofuran | ND | 10 | µg/L | | | | | | | U |
| Toluene | ND | 1.0 | µg/L | | | | | | | U |
| 1,2,3-Trichlorobenzene | 0.58 | 5.0 | µg/L | | | | | | | V-05, J |
| 1,2,4-Trichlorobenzene | 0.22 | 1.0 | µg/L | | | | | | | V-05, J |
| 1,3,5-Trichlorobenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,1,1-Trichloroethane | ND | 1.0 | µg/L | | | | | | | U |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032642 - SW-846 5030B

Blank (B032642-BLK1)

Prepared & Analyzed: 06/24/11

| | | | | | | | | | | |
|---|------|-----|------|------|--|------|--------|--|--|---|
| 1,1,2-Trichloroethane | ND | 1.0 | µg/L | | | | | | | U |
| Trichloroethylene | ND | 1.0 | µg/L | | | | | | | U |
| Trichlorofluoromethane (Freon 11) | ND | 2.0 | µg/L | | | | | | | U |
| 1,2,3-Trichloropropane | ND | 2.0 | µg/L | | | | | | | U |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | ND | 1.0 | µg/L | | | | | | | U |
| 1,2,4-Trimethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| 1,3,5-Trimethylbenzene | ND | 1.0 | µg/L | | | | | | | U |
| Vinyl Chloride | ND | 2.0 | µg/L | | | | | | | U |
| m+p Xylene | ND | 2.0 | µg/L | | | | | | | U |
| o-Xylene | ND | 1.0 | µg/L | | | | | | | U |
| Surrogate: 1,2-Dichloroethane-d4 | 24.6 | | µg/L | 25.0 | | 98.5 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.2 | | µg/L | 25.0 | | 97.0 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 24.8 | | µg/L | 25.0 | | 99.2 | 70-130 | | | |

LCS (B032642-BS1)

Prepared & Analyzed: 06/24/11

| | | | | | | | | | | |
|------------------------------------|------|------|------|------|--|------|--------|----|------|---|
| Acetone | 145 | 50 | µg/L | 100 | | 145 | 70-160 | 25 | | † |
| Acrylonitrile | 9.40 | 5.0 | µg/L | 10.0 | | 94.0 | 70-130 | 25 | | |
| tert-Amyl Methyl Ether (TAME) | 10.0 | 0.50 | µg/L | 10.0 | | 100 | 70-130 | 25 | | |
| Benzene | 10.0 | 1.0 | µg/L | 10.0 | | 100 | 70-130 | 25 | | |
| Bromobenzene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 25 | | |
| Bromochloromethane | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | 25 | | |
| Bromodichloromethane | 10.5 | 0.50 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |
| Bromoform | 8.74 | 1.0 | µg/L | 10.0 | | 87.4 | 70-130 | 25 | | |
| Bromomethane | 4.29 | 2.0 | µg/L | 10.0 | | 42.9 | 40-160 | 25 | | † |
| 2-Butanone (MEK) | 117 | 20 | µg/L | 100 | | 117 | 40-160 | 25 | | † |
| tert-Butyl Alcohol (TBA) | 104 | 20 | µg/L | 100 | | 104 | 40-160 | 25 | V-16 | † |
| n-Butylbenzene | 10.1 | 1.0 | µg/L | 10.0 | | 101 | 70-130 | 25 | | |
| sec-Butylbenzene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 25 | | |
| tert-Butylbenzene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 25 | | |
| tert-Butyl Ethyl Ether (TBEE) | 8.72 | 0.50 | µg/L | 10.0 | | 87.2 | 70-130 | 25 | | |
| Carbon Disulfide | 8.29 | 2.0 | µg/L | 10.0 | | 82.9 | 70-130 | 25 | | |
| Carbon Tetrachloride | 9.23 | 5.0 | µg/L | 10.0 | | 92.3 | 70-130 | 25 | | |
| Chlorobenzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |
| Chlorodibromomethane | 10.6 | 0.50 | µg/L | 10.0 | | 106 | 70-130 | 25 | | |
| Chloroethane | 7.94 | 2.0 | µg/L | 10.0 | | 79.4 | 70-130 | 25 | | |
| Chloroform | 10.1 | 2.0 | µg/L | 10.0 | | 101 | 70-130 | 25 | | |
| Chloromethane | 5.67 | 2.0 | µg/L | 10.0 | | 56.7 | 40-160 | 25 | | † |
| 2-Chlorotoluene | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 25 | | |
| 4-Chlorotoluene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 25 | | |
| 1,2-Dibromo-3-chloropropane (DBCP) | 9.02 | 5.0 | µg/L | 10.0 | | 90.2 | 70-130 | 25 | | |
| 1,2-Dibromoethane (EDB) | 10.8 | 0.50 | µg/L | 10.0 | | 108 | 70-130 | 25 | | |
| Dibromomethane | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | 25 | | |
| 1,2-Dichlorobenzene | 10.1 | 1.0 | µg/L | 10.0 | | 101 | 70-130 | 25 | | |
| 1,3-Dichlorobenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 25 | | |
| 1,4-Dichlorobenzene | 9.74 | 1.0 | µg/L | 10.0 | | 97.4 | 70-130 | 25 | | |
| trans-1,4-Dichloro-2-butene | 8.94 | 2.0 | µg/L | 10.0 | | 89.4 | 70-130 | 25 | | |
| Dichlorodifluoromethane (Freon 12) | 4.07 | 2.0 | µg/L | 10.0 | | 40.7 | 40-160 | 25 | | † |
| 1,1-Dichloroethane | 9.85 | 1.0 | µg/L | 10.0 | | 98.5 | 70-130 | 25 | | |
| 1,2-Dichloroethane | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 25 | | |
| 1,1-Dichloroethylene | 9.19 | 1.0 | µg/L | 10.0 | | 91.9 | 70-130 | 25 | | |
| cis-1,2-Dichloroethylene | 9.88 | 1.0 | µg/L | 10.0 | | 98.8 | 70-130 | 25 | | |
| trans-1,2-Dichloroethylene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 25 | | |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|--------------|
| Batch B032642 - SW-846 5030B | | | | | | | | | | |
| LCS (B032642-BS1) | | | | | | | | | | |
| Prepared & Analyzed: 06/24/11 | | | | | | | | | | |
| 1,2-Dichloropropane | 10.1 | 1.0 | µg/L | 10.0 | | 101 | 70-130 | | 25 | |
| 1,3-Dichloropropane | 10.5 | 0.50 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| 2,2-Dichloropropane | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 40-130 | | 25 | † |
| 1,1-Dichloropropene | 10.1 | 2.0 | µg/L | 10.0 | | 101 | 70-130 | | 25 | |
| cis-1,3-Dichloropropene | 9.24 | 0.50 | µg/L | 10.0 | | 92.4 | 70-130 | | 25 | |
| trans-1,3-Dichloropropene | 9.77 | 0.50 | µg/L | 10.0 | | 97.7 | 70-130 | | 25 | |
| Diethyl Ether | 9.87 | 2.0 | µg/L | 10.0 | | 98.7 | 70-130 | | 25 | |
| Diisopropyl Ether (DIPE) | 9.27 | 0.50 | µg/L | 10.0 | | 92.7 | 70-130 | | 25 | |
| 1,4-Dioxane | 117 | 50 | µg/L | 100 | | 117 | 40-130 | | 50 | V-16, V-20 † |
| Ethylbenzene | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | | 25 | |
| Hexachlorobutadiene | 10.4 | 0.50 | µg/L | 10.0 | | 104 | 70-130 | | 25 | |
| 2-Hexanone (MBK) | 118 | 10 | µg/L | 100 | | 118 | 70-160 | | 25 | † |
| Isopropylbenzene (Cumene) | 12.6 | 1.0 | µg/L | 10.0 | | 126 | 70-130 | | 25 | |
| p-Isopropyltoluene (p-Cymene) | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Methyl tert-Butyl Ether (MTBE) | 9.58 | 1.0 | µg/L | 10.0 | | 95.8 | 70-130 | | 25 | |
| Methylene Chloride | 9.59 | 5.0 | µg/L | 10.0 | | 95.9 | 70-130 | | 25 | |
| 4-Methyl-2-pentanone (MIBK) | 110 | 10 | µg/L | 100 | | 110 | 70-160 | | 25 | † |
| Naphthalene | 8.56 | 2.0 | µg/L | 10.0 | | 85.6 | 40-130 | | 25 | V-05 † |
| n-Propylbenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | | 25 | |
| Styrene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | | 25 | |
| 1,1,1,2-Tetrachloroethane | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | | 25 | |
| 1,1,2,2-Tetrachloroethane | 10.2 | 0.50 | µg/L | 10.0 | | 102 | 70-130 | | 25 | |
| Tetrachloroethylene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | | 25 | |
| Tetrahydrofuran | 8.79 | 10 | µg/L | 10.0 | | 87.9 | 70-130 | | 25 | J |
| Toluene | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | | 25 | |
| 1,2,3-Trichlorobenzene | 8.32 | 5.0 | µg/L | 10.0 | | 83.2 | 70-130 | | 25 | V-05 |
| 1,2,4-Trichlorobenzene | 8.43 | 1.0 | µg/L | 10.0 | | 84.3 | 70-130 | | 25 | V-05 |
| 1,3,5-Trichlorobenzene | 8.84 | 1.0 | µg/L | 10.0 | | 88.4 | 70-130 | | 25 | |
| 1,1,1-Trichloroethane | 10.7 | 1.0 | µg/L | 10.0 | | 107 | 70-130 | | 25 | |
| 1,1,2-Trichloroethane | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | | 25 | |
| Trichloroethylene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Trichlorofluoromethane (Freon 11) | 7.84 | 2.0 | µg/L | 10.0 | | 78.4 | 70-130 | | 25 | |
| 1,2,3-Trichloropropane | 9.17 | 2.0 | µg/L | 10.0 | | 91.7 | 70-130 | | 25 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 8.93 | 1.0 | µg/L | 10.0 | | 89.3 | 70-130 | | 25 | |
| 1,2,4-Trimethylbenzene | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | | 25 | |
| 1,3,5-Trimethylbenzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | | 25 | |
| Vinyl Chloride | 6.71 | 2.0 | µg/L | 10.0 | | 67.1 | 40-160 | | 25 | † |
| m+p Xylene | 22.1 | 2.0 | µg/L | 20.0 | | 110 | 70-130 | | 25 | |
| o-Xylene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | | 25 | |
| Surrogate: 1,2-Dichloroethane-d4 | 23.8 | | µg/L | 25.0 | | 95.0 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.8 | | µg/L | 25.0 | | 99.1 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 25.1 | | µg/L | 25.0 | | 101 | 70-130 | | | |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|
|---------|--------|-----------------|-------|-------------|---------------|------|-------------|-----|-----------|-------|

Batch B032642 - SW-846 5030B

LCS Dup (B032642-BSD1)

Prepared & Analyzed: 06/24/11

| | | | | | | | | | | |
|---|------|------|------|------|--|---------------|--------|-------|----|----------------|
| Acetone | 136 | 50 | µg/L | 100 | | 136 | 70-160 | 6.58 | 25 | † |
| Acrylonitrile | 9.21 | 5.0 | µg/L | 10.0 | | 92.1 | 70-130 | 2.04 | 25 | |
| tert-Amyl Methyl Ether (TAME) | 10.2 | 0.50 | µg/L | 10.0 | | 102 | 70-130 | 1.19 | 25 | |
| Benzene | 9.79 | 1.0 | µg/L | 10.0 | | 97.9 | 70-130 | 2.32 | 25 | |
| Bromobenzene | 10.0 | 1.0 | µg/L | 10.0 | | 100 | 70-130 | 2.47 | 25 | |
| Bromochloromethane | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 2.27 | 25 | |
| Bromodichloromethane | 10.5 | 0.50 | µg/L | 10.0 | | 105 | 70-130 | 0.476 | 25 | |
| Bromoform | 8.99 | 1.0 | µg/L | 10.0 | | 89.9 | 70-130 | 2.82 | 25 | |
| Bromomethane | 4.97 | 2.0 | µg/L | 10.0 | | 49.7 | 40-160 | 14.7 | 25 | † |
| 2-Butanone (MEK) | 111 | 20 | µg/L | 100 | | 111 | 40-160 | 5.14 | 25 | † |
| tert-Butyl Alcohol (TBA) | 98.4 | 20 | µg/L | 100 | | 98.4 | 40-160 | 5.38 | 25 | V-16 † |
| n-Butylbenzene | 9.96 | 1.0 | µg/L | 10.0 | | 99.6 | 70-130 | 1.79 | 25 | |
| sec-Butylbenzene | 10.0 | 1.0 | µg/L | 10.0 | | 100 | 70-130 | 1.38 | 25 | |
| tert-Butylbenzene | 10.1 | 1.0 | µg/L | 10.0 | | 101 | 70-130 | 1.08 | 25 | |
| tert-Butyl Ethyl Ether (TBEE) | 8.73 | 0.50 | µg/L | 10.0 | | 87.3 | 70-130 | 0.115 | 25 | |
| Carbon Disulfide | 7.98 | 2.0 | µg/L | 10.0 | | 79.8 | 70-130 | 3.81 | 25 | |
| Carbon Tetrachloride | 8.71 | 5.0 | µg/L | 10.0 | | 87.1 | 70-130 | 5.80 | 25 | |
| Chlorobenzene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 0.477 | 25 | |
| Chlorodibromomethane | 10.8 | 0.50 | µg/L | 10.0 | | 108 | 70-130 | 2.15 | 25 | |
| Chloroethane | 7.68 | 2.0 | µg/L | 10.0 | | 76.8 | 70-130 | 3.33 | 25 | |
| Chloroform | 9.88 | 2.0 | µg/L | 10.0 | | 98.8 | 70-130 | 2.50 | 25 | |
| Chloromethane | 5.42 | 2.0 | µg/L | 10.0 | | 54.2 | 40-160 | 4.51 | 25 | † |
| 2-Chlorotoluene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 2.48 | 25 | |
| 4-Chlorotoluene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | 0.365 | 25 | |
| 1,2-Dibromo-3-chloropropane (DBCP) | 9.01 | 5.0 | µg/L | 10.0 | | 90.1 | 70-130 | 0.111 | 25 | |
| 1,2-Dibromoethane (EDB) | 10.7 | 0.50 | µg/L | 10.0 | | 107 | 70-130 | 0.372 | 25 | |
| Dibromomethane | 10.8 | 1.0 | µg/L | 10.0 | | 108 | 70-130 | 1.75 | 25 | |
| 1,2-Dichlorobenzene | 9.99 | 1.0 | µg/L | 10.0 | | 99.9 | 70-130 | 1.49 | 25 | |
| 1,3-Dichlorobenzene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 1.46 | 25 | |
| 1,4-Dichlorobenzene | 9.76 | 1.0 | µg/L | 10.0 | | 97.6 | 70-130 | 0.205 | 25 | |
| trans-1,4-Dichloro-2-butene | 8.39 | 2.0 | µg/L | 10.0 | | 83.9 | 70-130 | 6.35 | 25 | |
| Dichlorodifluoromethane (Freon 12) | 3.66 | 2.0 | µg/L | 10.0 | | 36.6 * | 40-160 | 10.6 | 25 | L-07 † |
| 1,1-Dichloroethane | 9.68 | 1.0 | µg/L | 10.0 | | 96.8 | 70-130 | 1.74 | 25 | |
| 1,2-Dichloroethane | 10.6 | 1.0 | µg/L | 10.0 | | 106 | 70-130 | 2.19 | 25 | |
| 1,1-Dichloroethylene | 9.12 | 1.0 | µg/L | 10.0 | | 91.2 | 70-130 | 0.765 | 25 | |
| cis-1,2-Dichloroethylene | 9.57 | 1.0 | µg/L | 10.0 | | 95.7 | 70-130 | 3.19 | 25 | |
| trans-1,2-Dichloroethylene | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 70-130 | 2.80 | 25 | |
| 1,2-Dichloropropane | 10.1 | 1.0 | µg/L | 10.0 | | 101 | 70-130 | 0.297 | 25 | |
| 1,3-Dichloropropane | 10.2 | 0.50 | µg/L | 10.0 | | 102 | 70-130 | 2.22 | 25 | |
| 2,2-Dichloropropane | 10.2 | 1.0 | µg/L | 10.0 | | 102 | 40-130 | 7.20 | 25 | † |
| 1,1-Dichloropropene | 9.73 | 2.0 | µg/L | 10.0 | | 97.3 | 70-130 | 3.43 | 25 | |
| cis-1,3-Dichloropropene | 8.96 | 0.50 | µg/L | 10.0 | | 89.6 | 70-130 | 3.08 | 25 | |
| trans-1,3-Dichloropropene | 9.69 | 0.50 | µg/L | 10.0 | | 96.9 | 70-130 | 0.822 | 25 | |
| Diethyl Ether | 10.2 | 2.0 | µg/L | 10.0 | | 102 | 70-130 | 3.09 | 25 | |
| Diisopropyl Ether (DIPE) | 9.42 | 0.50 | µg/L | 10.0 | | 94.2 | 70-130 | 1.61 | 25 | |
| 1,4-Dioxane | 114 | 50 | µg/L | 100 | | 114 | 40-130 | 2.74 | 50 | V-16, V-20 † ‡ |
| Ethylbenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 2.46 | 25 | |
| Hexachlorobutadiene | 10.1 | 0.50 | µg/L | 10.0 | | 101 | 70-130 | 2.92 | 25 | |
| 2-Hexanone (MBK) | 111 | 10 | µg/L | 100 | | 111 | 70-160 | 6.12 | 25 | † |
| Isopropylbenzene (Cumene) | 12.2 | 1.0 | µg/L | 10.0 | | 122 | 70-130 | 2.98 | 25 | |
| p-Isopropyltoluene (p-Cymene) | 10.0 | 1.0 | µg/L | 10.0 | | 100 | 70-130 | 4.97 | 25 | |
| Methyl tert-Butyl Ether (MTBE) | 9.98 | 1.0 | µg/L | 10.0 | | 99.8 | 70-130 | 4.09 | 25 | |

QUALITY CONTROL

Volatile Organic Compounds by GC/MS - Quality Control

| Analyte | Result | Reporting Limit | Units | Spike Level | Source Result | %REC | %REC Limits | RPD | RPD Limit | Notes |
|---|--------|-----------------|-------|-------------|---------------|------|-------------|-------|-----------|--------|
| Batch B032642 - SW-846 5030B | | | | | | | | | | |
| LCS Dup (B032642-BSD1) | | | | | | | | | | |
| Prepared & Analyzed: 06/24/11 | | | | | | | | | | |
| Methylene Chloride | 9.66 | 5.0 | µg/L | 10.0 | | 96.6 | 70-130 | 0.727 | 25 | |
| 4-Methyl-2-pentanone (MIBK) | 107 | 10 | µg/L | 100 | | 107 | 70-160 | 2.55 | 25 | † |
| Naphthalene | 8.31 | 2.0 | µg/L | 10.0 | | 83.1 | 40-130 | 2.96 | 25 | V-05 † |
| n-Propylbenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 0.482 | 25 | |
| Styrene | 11.0 | 1.0 | µg/L | 10.0 | | 110 | 70-130 | 0.911 | 25 | |
| 1,1,1,2-Tetrachloroethane | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 2.04 | 25 | |
| 1,1,2,2-Tetrachloroethane | 10.1 | 0.50 | µg/L | 10.0 | | 101 | 70-130 | 0.790 | 25 | |
| Tetrachloroethylene | 10.5 | 1.0 | µg/L | 10.0 | | 105 | 70-130 | 4.45 | 25 | |
| Tetrahydrofuran | 9.13 | 10 | µg/L | 10.0 | | 91.3 | 70-130 | 3.79 | 25 | J |
| Toluene | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | 2.49 | 25 | |
| 1,2,3-Trichlorobenzene | 7.93 | 5.0 | µg/L | 10.0 | | 79.3 | 70-130 | 4.80 | 25 | V-05 |
| 1,2,4-Trichlorobenzene | 8.01 | 1.0 | µg/L | 10.0 | | 80.1 | 70-130 | 5.11 | 25 | V-05 |
| 1,3,5-Trichlorobenzene | 8.64 | 1.0 | µg/L | 10.0 | | 86.4 | 70-130 | 2.29 | 25 | |
| 1,1,1-Trichloroethane | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | 3.53 | 25 | |
| 1,1,2-Trichloroethane | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 0.289 | 25 | |
| Trichloroethylene | 10.3 | 1.0 | µg/L | 10.0 | | 103 | 70-130 | 2.21 | 25 | |
| Trichlorofluoromethane (Freon 11) | 7.52 | 2.0 | µg/L | 10.0 | | 75.2 | 70-130 | 4.17 | 25 | |
| 1,2,3-Trichloropropane | 9.26 | 2.0 | µg/L | 10.0 | | 92.6 | 70-130 | 0.977 | 25 | |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | 8.91 | 1.0 | µg/L | 10.0 | | 89.1 | 70-130 | 0.224 | 25 | |
| 1,2,4-Trimethylbenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 1.26 | 25 | |
| 1,3,5-Trimethylbenzene | 10.4 | 1.0 | µg/L | 10.0 | | 104 | 70-130 | 1.15 | 25 | |
| Vinyl Chloride | 6.67 | 2.0 | µg/L | 10.0 | | 66.7 | 40-160 | 0.598 | 25 | † |
| m+p Xylene | 21.6 | 2.0 | µg/L | 20.0 | | 108 | 70-130 | 2.43 | 25 | |
| o-Xylene | 10.9 | 1.0 | µg/L | 10.0 | | 109 | 70-130 | 0.910 | 25 | |
| Surrogate: 1,2-Dichloroethane-d4 | 23.8 | | µg/L | 25.0 | | 95.2 | 70-130 | | | |
| Surrogate: Toluene-d8 | 24.4 | | µg/L | 25.0 | | 97.6 | 70-130 | | | |
| Surrogate: 4-Bromofluorobenzene | 25.3 | | µg/L | 25.0 | | 101 | 70-130 | | | |

FLAG/QUALIFIER SUMMARY

- * QC result is outside of established limits.
 - † Wide recovery limits established for difficult compound.
 - ‡ Wide RPD limits established for difficult compound.
 - # Data exceeded client recommended or regulatory level
- Percent recoveries and relative percent differences (RPDs) are determined by the software using values in the calculation which have not been rounded.
- J Detected but below the Reporting Limit (lowest calibration standard); therefore, result is an estimated concentration (CLP J-Flag).
 - L-02 Laboratory fortified blank/laboratory control sample recovery and duplicate recoveries outside of control limits. Data validation is not affected since all results are "not detected" for associated samples in this batch and bias is on the high side.
 - L-07 Either laboratory fortified blank/laboratory control sample or duplicate recovery is outside of control limits, but the other is within limits. RPD between the two LFB/LCS results is within method specified criteria.
 - U Analyte included in the analysis, but not detected
 - V-05 Continuing calibration did not meet method specifications and was biased on the low side for this compound. Increased uncertainty is associated with the reported value which is likely to be biased on the low side.
 - V-16 Response factor is less than method specified minimum acceptable value. Reduced precision and accuracy are associated with reported result.
 - V-20 Continuing calibration did not meet method specifications and was biased on the high side. Data validation is not affected since sample result was "not detected" for this compound.

CERTIFICATIONS

Certified Analyses included in this Report

| Analyte | Certifications |
|------------------------------------|-------------------|
| <i>SW-846 8260C in Water</i> | |
| Acetone | CT,NH,NY,NC,ME |
| Acrylonitrile | CT,NY,NC,RI,ME |
| tert-Amyl Methyl Ether (TAME) | NH,NY,NC,ME |
| Benzene | CT,NH,NY,NC,RI,ME |
| Bromobenzene | NC |
| Bromochloromethane | NH,NY,NC,ME |
| Bromodichloromethane | CT,NH,NY,NC,RI,ME |
| Bromoform | CT,NH,NY,NC,RI,ME |
| Bromomethane | CT,NH,NY,NC,RI,ME |
| 2-Butanone (MEK) | CT,NH,NY,NC,ME |
| tert-Butyl Alcohol (TBA) | NH,NY,NC,ME |
| n-Butylbenzene | NY,NC,ME |
| sec-Butylbenzene | NY,NC,ME |
| tert-Butylbenzene | NY,NC,ME |
| tert-Butyl Ethyl Ether (TBEE) | NH,NY,NC,ME |
| Carbon Disulfide | CT,NH,NY,NC,ME |
| Carbon Tetrachloride | CT,NH,NY,NC,RI,ME |
| Chlorobenzene | CT,NH,NY,NC,RI,ME |
| Chlorodibromomethane | CT,NH,NY,NC,RI,ME |
| Chloroethane | CT,NH,NY,NC,RI,ME |
| Chloroform | CT,NH,NY,NC,RI,ME |
| Chloromethane | CT,NH,NY,NC,RI,ME |
| 2-Chlorotoluene | NY,NC,ME |
| 4-Chlorotoluene | NY,NC,ME |
| 1,2-Dibromo-3-chloropropane (DBCP) | NC |
| 1,2-Dibromoethane (EDB) | NC |
| Dibromomethane | NH,NY,NC,ME |
| 1,2-Dichlorobenzene | CT,NY,NC,RI,ME |
| 1,3-Dichlorobenzene | CT,NH,NY,NC,RI,ME |
| 1,4-Dichlorobenzene | CT,NH,NY,NC,RI,ME |
| trans-1,4-Dichloro-2-butene | NH,NY,NC,ME |
| Dichlorodifluoromethane (Freon 12) | NH,NY,NC,RI,ME |
| 1,1-Dichloroethane | CT,NH,NY,NC,RI,ME |
| 1,2-Dichloroethane | CT,NH,NY,NC,RI,ME |
| 1,1-Dichloroethylene | CT,NH,NY,NC,RI,ME |
| cis-1,2-Dichloroethylene | NC,ME |
| trans-1,2-Dichloroethylene | CT,NH,NY,NC,RI,ME |
| 1,2-Dichloropropane | CT,NH,NY,NC,RI,ME |
| 1,3-Dichloropropane | NY,NC,ME |
| 2,2-Dichloropropane | NH,NY,NC,ME |
| 1,1-Dichloropropene | NH,NY,NC,ME |
| cis-1,3-Dichloropropene | CT,NH,NY,NC,RI,ME |
| trans-1,3-Dichloropropene | CT,NH,NY,NC,RI,ME |
| Diethyl Ether | NC |
| Diisopropyl Ether (DIPE) | NH,NY,NC,ME |
| 1,4-Dioxane | NC |
| Ethylbenzene | CT,NH,NY,NC,RI,ME |

CERTIFICATIONS

Certified Analyses included in this Report

| Analyte | Certifications |
|---|-------------------|
| <i>SW-846 8260C in Water</i> | |
| Hexachlorobutadiene | CT,NH,NY,NC,ME |
| 2-Hexanone (MBK) | CT,NH,NY,NC,ME |
| Isopropylbenzene (Cumene) | NY,NC,ME |
| p-Isopropyltoluene (p-Cymene) | CT,NH,NY,NC,ME |
| Methyl tert-Butyl Ether (MTBE) | CT,NH,NY,NC,ME |
| Methylene Chloride | CT,NH,NY,NC,RI,ME |
| 4-Methyl-2-pentanone (MIBK) | CT,NH,NY,NC,ME |
| Naphthalene | NH,NY,NC,ME |
| n-Propylbenzene | CT,NH,NY,NC,ME |
| Styrene | CT,NH,NY,NC,ME |
| 1,1,1,2-Tetrachloroethane | CT,NH,NY,NC,ME |
| 1,1,2,2-Tetrachloroethane | CT,NH,NY,NC,RI,ME |
| Tetrachloroethylene | CT,NH,NY,NC,RI,ME |
| Tetrahydrofuran | NC |
| Toluene | CT,NH,NY,NC,RI,ME |
| 1,2,3-Trichlorobenzene | NH,NY,NC,ME |
| 1,2,4-Trichlorobenzene | CT,NH,NY,NC,ME |
| 1,3,5-Trichlorobenzene | NC,ME |
| 1,1,1-Trichloroethane | CT,NH,NY,NC,RI,ME |
| 1,1,2-Trichloroethane | CT,NH,NY,NC,RI,ME |
| Trichloroethylene | CT,NH,NY,NC,RI,ME |
| Trichlorofluoromethane (Freon 11) | CT,NH,NY,NC,RI,ME |
| 1,2,3-Trichloropropane | NH,NY,NC,ME |
| 1,1,2-Trichloro-1,2,2-trifluoroethane (Freon 113) | NC |
| 1,2,4-Trimethylbenzene | NY,NC,ME |
| 1,3,5-Trimethylbenzene | NY,NC,ME |
| Vinyl Chloride | CT,NH,NY,NC,RI,ME |
| m+p Xylene | CT,NH,NY,NC,RI,ME |
| o-Xylene | CT,NH,NY,NC,RI,ME |

The CON-TEST Environmental Laboratory operates under the following certifications and accreditations:

| Code | Description | Number | Expires |
|------|--|---------------|------------|
| AIHA | American Industrial Hygiene Association | 100033 | 01/1/2012 |
| MA | Massachusetts DEP | M-MA100 | 06/30/2011 |
| CT | Connecticut Department of Public Health | PH-0567 | 09/30/2011 |
| NY | New York State Department of Health | 10899 NELAP | 04/1/2012 |
| NH | New Hampshire Environmental Lab | 2516 NELAP | 02/5/2012 |
| RI | Rhode Island Department of Health | LAO00112 | 12/30/2011 |
| NC | North Carolina Div. of Water Quality | 652 | 12/31/2011 |
| NJ | New Jersey DEP | MA007 NELAP | 06/30/2012 |
| FL | Florida Department of Health | E871027 NELAP | 06/30/2011 |
| VT | Vermont Department of Health Lead Laboratory | LL015036 | 07/30/2011 |
| WA | State of Washington Department of Ecology | C2065 | 02/23/2012 |
| ME | State of Maine | 2011028 | 06/9/2013 |



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 www.contestlabs.com

CHAIN OF CUSTODY RECORD
 11F075C

39 SPRUCE ST, 2ND FLOOR
 EAST LONGMEADOW, MA 01028

Company Name: WFB Inc.
 Address: 57 Hill Rd
Concord NH 03301
 Attention: S. Donaldson
 Project Location: 4805 Main St, Wartsfield, VT
 Sampled By: S. Donaldson

| | | | |
|---|---|----------------------------|--|
| Telephone: <u>(603) 364-4190</u> | Project # <u>4400001</u> | Client PO # <u>6000001</u> | # of containers **Preserved |
| DATA DELIVERY (check one): <input type="checkbox"/> FAX <input checked="" type="checkbox"/> EMAIL <input checked="" type="checkbox"/> WEBSITE CLIENT | <input type="checkbox"/> EXCEL <input checked="" type="checkbox"/> PDF <input type="checkbox"/> GIS KEY | | ~Cont. Code: A=amber glass G=glass P=plastic ST=sterile V=vial S=summary can T=tealbar bag O=Other |

Proposal Provided? (For Billing purposes) yes no
 State Form Required? yes no

| Field ID | Sample Description | Lab # | Start Date/Time | Stop Date/Time | Comp-site | Grab | *Matrix Code | Conc. Code | ANALYSIS REQUESTED | Client Comments: |
|----------|--------------------|-------|-----------------|----------------|-----------|------|--------------|------------|--------------------|------------------|
| -01 | MW-1 | | 6/16/11 10:00 | | | X | GW | U | | |
| -02 | MW-2 | | 1:15 | | | X | GW | U | | |
| -03 | MW-4 | | 10:35 | | | X | GW | U | | |

Please use the following codes to let Con-Test know if a specific sample may be high in concentration in Matrix/Conc. Code Box:
 H - High; M - Medium; L - Low; C - Clean; U - Unknown

Relinquished by: (signature) [Signature] Date/Time: 6/20/11 14:00

Received by: (signature) [Signature] Date/Time: 6-21-11 11:00

Relinquished by: (signature) [Signature] Date/Time: 6-21-11 17:40

Received by: (signature) [Signature] Date/Time: 6/21/11 17:40

Turnaround **
 7-Day
 10-Day
 Other RUSH
 *24-Hr *48-Hr
 *72-Hr *4-Day
 * Require lab approval

Detection Limit Requirements
 Regulations? _____
 Data Enhancement Project/RCP? Y N
 Special Requirements or DLS: _____

**Matrix Code:
 GW = groundwater
 WW = wastewater
 DW = drinking water
 A = air
 S = soil/solid
 SL = sludge
 O = other

**Preservation Codes:
 I = Iced
 H = HCL
 M = Methanol
 N = Nitric Acid
 S = Sulfuric Acid
 B = Sodium bisulfate
 O = Other
 X = Na hydroxide
 T = Na thiosulfate

** TURNAROUND TIME STARTS AT 9:00 A.M. THE DAY AFTER SAMPLE RECEIPT UNLESS THERE ARE QUESTIONS ON YOUR CHAIN. IF THIS FORM IS NOT FILLED OUT COMPLETELY OR IS INCORRECT, TURNAROUND TIME WILL NOT START UNTIL ALL QUESTIONS ARE ANSWERED BY OUR CLIENT.

AIHA, NELAC & WBE/DBE Certified

39 Spruce St.
 East Longmeadow, MA. 01028
 P: 413-525-2332
 F: 413-525-6405
 www.contestlabs.com



Sample Receipt Checklist

CLIENT NAME: WFB Inc. RECEIVED BY: [Signature] DATE: 02/11

- 1) Was the chain(s) of custody relinquished and signed? Yes No No CoC Included
- 2) Does the chain agree with the samples? Yes No
 If not, explain: _____
- 3) Are all the samples in good condition? Yes No
 If not, explain: _____

4) How were the samples received:
 On Ice Direct from Sampling Ambient In Cooler(s)
 Were the samples received in Temperature Compliance of (2-6°C)? Yes No N/A
 Temperature °C by Temp blank _____ Temperature °C by Temp gun 2.7°C

5) Are there Dissolved samples for the lab to filter? Yes No
 Who was notified _____ Date _____ Time _____

6) Are there any RUSH or SHORT HOLDING TIME samples? Yes No
 Who was notified _____ Date _____ Time _____

7) Location where samples are stored: 19

Permission to subcontract samples? Yes No
 (Walk-in clients only) if not already approved
 Client Signature: _____

Containers received at Con-Test

| | # of containers | | # of containers |
|--------------------------------|-----------------|-----------------------|-----------------|
| 1 Liter Amber | | 8 oz amber/clear jar | |
| 500 mL Amber | | 4 oz amber/clear jar | |
| 250 mL Amber (8oz amber) | | 2 oz amber/clear jar | |
| 1 Liter Plastic | | Air Cassette | |
| 500 mL Plastic | | Hg/Hopcalite Tube | |
| 250 mL plastic | | Plastic Bag / Ziploc | |
| 40 mL Vial - type listed below | 9 | PM 2.5 / PM 10 | |
| Colisure / bacteria bottle | | PUF Cartridge | |
| Dissolved Oxygen bottle | | SOC Kit | |
| Encore | | TO-17 Tubes | |
| Flashpoint bottle | | Non-ConTest Container | |
| Perchlorate Kit | | Other glass jar | |
| Other | | Other | |

Laboratory Comments: _____

40 mL vials: # HCl 9 # Methanol _____
 # Bisulfate _____ # DI Water _____
 # Thiosulfate _____ Unpreserved _____

Time and Date Frozen: _____

Do all samples have the proper Acid pH: Yes No N/A _____
 Do all samples have the proper Base pH: Yes No N/A _____