

**Groundwater Monitoring Report
Semi-Annual Sampling Event
First Half of 2020**

**Capital Center
705 Broadway
Albany, New York
BCP Site No. C401070**

CHA Project Number: 031647.000

Prepared for:

*FC 705 Broadway LLC
c/o Pioneer Companies
333 West Washington Street, Suite 600
Syracuse, New York 13202*

Prepared by:

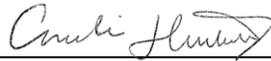


*III Winners Circle
Albany, New York 12205
Phone: (518) 453-4500
Fax: (518) 453-4773*

July 14, 2020

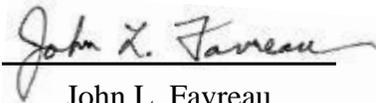
This report has been prepared and reviewed by the following qualified personnel employed by CHA.

Report Prepared By:



Caroline Hurlburt
Scientist I

Report Reviewed By:



John L. Favreau
Senior Scientist V

TABLE OF CONTENTS

1.0 Introduction..... 1
2.0 Sampling Procedures 2
3.0 Groundwater Flow Direction 3
4.0 Laboratory Analytical Results 3
5.0 Future Activities..... 4

FIGURES

Figure 1 – Site Location Map

Figure 2 – Site Plan

TABLES

Table 1 – Groundwater Elevations

Table 2 – Groundwater Analytical Data

APPENDIX

Appendix A – Monitoring Well Sampling Log

Appendix B – Laboratory Analytical Report

1.0 INTRODUCTION

This report has been prepared by CHA Consulting, Inc. (CHA) on behalf of FC 705 Broadway LLC and Pioneer Companies to describe the field activities and present the findings associated with the first of two semiannual groundwater monitoring events in 2020 at the Capital Center Site (the Site), located at 705 Broadway in the City of Albany, Albany County, New York (Figure 1). The semiannual monitoring event conducted during the first half of the year was completed by CHA on April 9, 2020. The monitoring event was conducted in accordance with the New York State Department of Environmental Conservation (NYSDEC)-approved Site Management Plan (SMP), prepared by CHA in December 2017. The SMP includes periodic groundwater monitoring as a means to evaluate the effectiveness of the implemented remedy (soil excavation) and natural attenuation.

Six groundwater monitoring wells (MW-1 to MW-6) were installed across the Site in March 2018, all to a depth of 25 feet below ground surface (bgs). Monitoring well construction logs are included in Appendix C of the SMP. Monitoring wells MW-3 through MW-6 were decommissioned in October 2018, based on consistent quarterly analytical results from these wells showing concentrations of volatile organic compounds (VOCs) and semivolatile organic compounds (SVOCs) below the TOGS 1.1.1 standard values. Monitoring well MW-2 was subsequently decommissioned in January 2020, based on analytical results from this well showing VOC and SVOC concentrations (particularly MTBE) below the TOGS 1.1.1 standard values. Monitoring well MW-1, located downgradient of the former on-site gasoline station building, was left in place to facilitate continued monitoring in this area; however, the PVC well riser was extended by five feet to accommodate filling and grading of the Site with crushed stone for construction purposes. The riser was subsequently extended another five feet as filling and grading progressed in this area of the Site. . A site plan showing the locations of the existing and former monitoring wells is included as Figure 2.

2.0 SAMPLING PROCEDURES

On April 9, 2020, CHA personnel conducted groundwater monitoring activities at well MW-1 in accordance with the NYSDEC-approved SMP and CHA's standard operating procedure (SOP) for monitoring well sampling.

Prior to conducting groundwater sampling activities, the water level in the well was measured from the top of the well riser using an electronic water level meter. The well was then purged via low-flow sampling methods, using a portable bladder pump and dedicated tubing. During purging, the following parameters were measured at approximately five-minute intervals using a YSI water quality meter equipped with a flow-through cell: specific conductivity; temperature; pH; and oxidation/reduction potential (ORP). Turbidity was also measured using a separate turbidity meter. Field parameter measurements and observations were recorded and are documented in the Monitoring Well Sampling Log included as Appendix A.

Following stabilization of field parameters, a groundwater sample was collected from the well via the portable bladder pump and dedicated tubing. The sample was collected in laboratory-provided, pre-preserved containers, which were labeled and then placed in a rigid cooler with ice, pending submittal to the laboratory.

Following sampling activities, the groundwater sample was delivered by CHA to Alpha Analytical's (Alpha) Service Center in Albany, New York, for subsequent transport by Alpha to its laboratory in Westborough, Massachusetts, in accordance with proper chain-of-custody protocol. Alpha is currently certified by the New York State Department of Health's (NYSDOH) Environmental Laboratory Approval Program (ELAP). The groundwater sample was analyzed for VOCs and SVOCs via EPA Methods 8260 and 8270, respectively, in accordance with applicable NYSDOH QA/QC procedures. A trip blank sample, prepared by the laboratory, accompanied the groundwater sample to the laboratory and was to have been analyzed for VOCs only; however, the sample identification was inadvertently omitted from the chain-of-custody and therefore was not analyzed. Groundwater samples from the Site were historically also analyzed for metals; however, based on evaluation of detected metals during the past few monitoring events, the NYSDEC approved the discontinuation of metals analysis, beginning with the April 2020 monitoring event.

3.0 GROUNDWATER FLOW DIRECTION

Groundwater elevation data from monitoring well MW-1 is included in Table 1. It should be noted that water level measurements are required from a minimum of three locations to triangulate and determine groundwater flow direction. Due to availability of only one data point, a groundwater contour map was not generated; however, the water level measurement at well MW-1 indicates a relative groundwater elevation similar to that exhibited during past monitoring events at this location. Based on this data, along with historical data and local topography, the expected groundwater flow direction continues to be generally to the east, toward the Hudson River.

4.0 LABORATORY ANALYTICAL RESULTS

Groundwater analytical results were compared to the groundwater standards and/or guidance values listed in the NYSDEC Division of Water Technical and Operational Guidance Series 1.1.1: Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations (TOGS 1.1.1). Those parameters in excess of their respective standards or guidance values are shaded in Table 2. Appendix B includes the laboratory deliverables package for the April 2020 sampling event. A discussion of the analytical results for the groundwater samples collected from each monitoring well is provided below.

MW-1

In the sample collected from well MW-1, the VOC methyl tert-butyl ether (MTBE) was detected in exceedance of its TOGS 1.1.1 standard; however, the detected concentration was significantly reduced (nearly 80 percent) as compared to the August 2019 sampling results. Acetone was also detected in the sample, but at a concentration below its TOGS 1.1.1 standard. It should be noted that acetone is a common laboratory artifact and as a result, the detection of acetone in the MW-1 sample may be a result of laboratory contamination. No SVOCs were detected at concentrations above laboratory reporting limits.

Summary

The VOC MTBE was detected at the location of well MW-1 (in the area downgradient of the former on-site gasoline station) at a concentration exceeding its TOGS 1.1.1 standard; however, the detected concentration (34 µg/l) represents a significant decrease from the previous sampling event in August 2019 and indicates a continued decreasing trend in the concentration of MTBE at this location. The continued decreasing trend in MTBE concentrations at well MW-1 is indicative

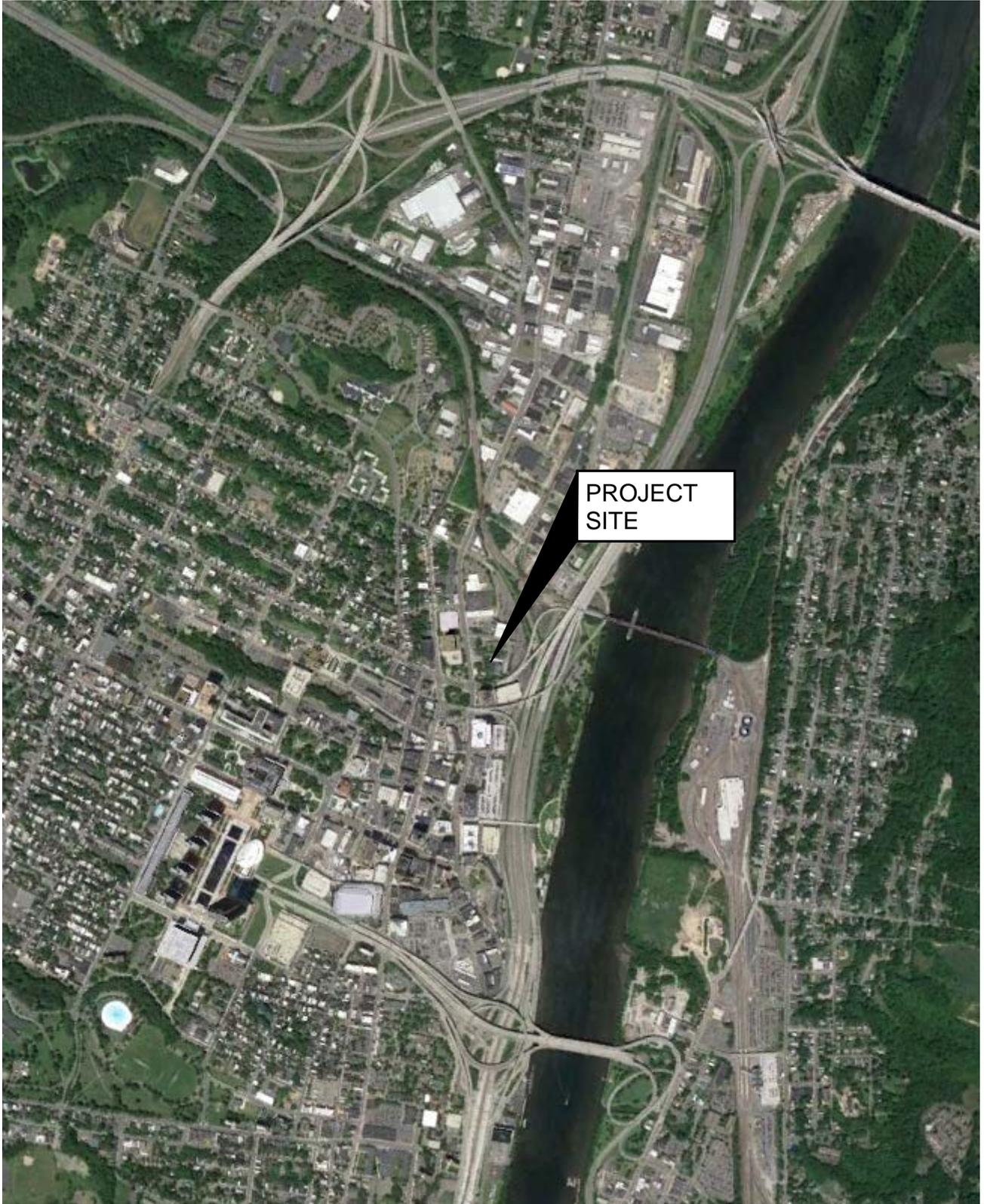
of the effectiveness of the implemented remedial activities at the Site, along with natural attenuation.

5.0 FUTURE ACTIVITIES

The next semiannual groundwater monitoring event is scheduled to occur in October 2020. Given the recent decreasing trend in MTBE concentration at well MW-1 and in particular, the significant decrease exhibited during the April 2020 monitoring event, if the analytical results from the October 2020 monitoring event indicate a continued decreasing trend, it is anticipated that CHA will recommend that groundwater monitoring be discontinued and monitoring well MW-1 be properly decommissioned.

FIGURES

SOURCE: Google Earth



V:\Projects\NY\K4131647\Reports\Site Management Plan\Figures\DECU\updated_Figure1_KE.doc

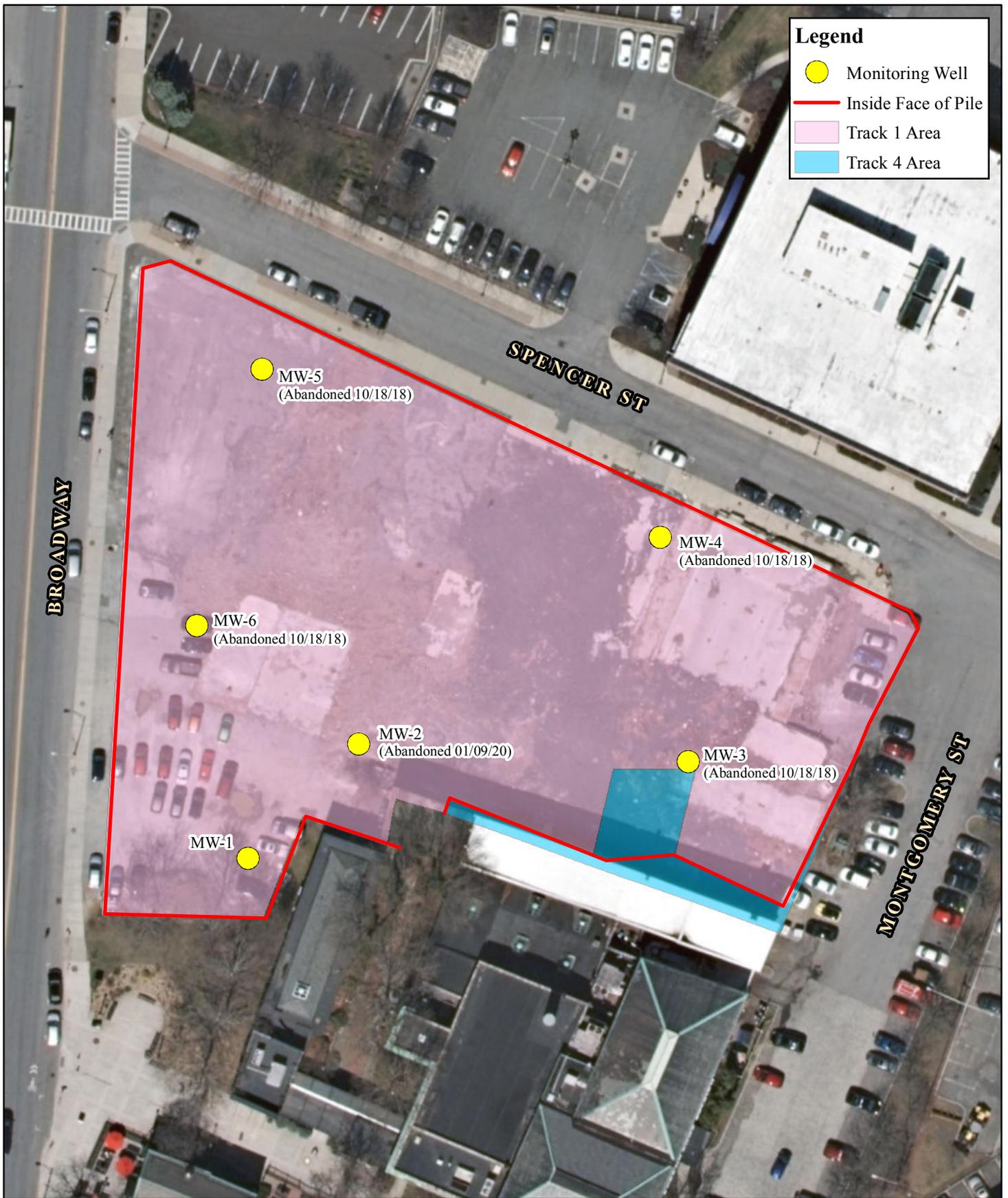


111 Winners Circle, P.O. Box 6289 • Albany, NY 12206-0289
Main: (518)463-4600 • www.oloughharbour.com

NOT TO SCALE

DATE: December 2017

FIGURE 1
SITE LOCATION MAP
CAPITAL CENTER PROJECT
ALBANY COUNTY, NEW YORK



Legend

- Monitoring Well
- Inside Face of Pile
- Track 1 Area
- Track 4 Area

| | | |
|---|---|--|
|  |  | <p>Figure 2 - Site Plan</p> <p>Capital Center Site Quackenbush Square, 705 Broadway City of Albany, Albany County, New York</p> |
| <p>Scale 1" = 60'</p> | <p>CHA Project No. 31647</p> | <p><i>Service Layer Credits: Aerial imagery courtesy of Esri, DigitalGlobe, GeoEye, Earthstar Image Date: 6/7/2015</i></p> |

TABLES

**Table 1
Groundwater Elevations**

| Well ID | MW-1 | MW-2 | MW-3 | MW-4 | MW-5 | MW-6 |
|---------------------|--------|----------------|----------------|----------------|----------------|----------------|
| TOR Elevations (ft) | 103.27 | 101.46 | 96.23 | 94.22 | 101.27 | 101.88 |
| 3/27/2018 | 98.93 | 96.78 | 90.53 | 89.42 | 100.02 | 96.21 |
| 6/6/2018 | 98.42 | 96.17 | 89.98 | 89.61 | 99.52 | 99.21 |
| 9/27/2018 | 98.46 | 96.70 | 85.83 | 86.27 | 99.99 | 99.46 |
| 12/19/2018 | 97.27 | 96.20 | decommissioned | decommissioned | decommissioned | decommissioned |
| 3/20/2019 | 99.11 | 97.15 | decommissioned | decommissioned | decommissioned | decommissioned |
| 8/23/2019 | 98.73 | 97.84 | decommissioned | decommissioned | decommissioned | decommissioned |
| 4/9/2020 | 98.92 | decommissioned | decommissioned | decommissioned | decommissioned | decommissioned |

Note: MW-3 through MW-6 decommissioned and removed in October 2018
 MW-2 decommissioned and grouted in place in January 2020
 TOR Elevation corrected to 113.27 for 4/9/2020 event, due to raising of well riser following filling/grading of site

Table 2
Capital Center, Albany, NY
Groundwater Analytical Monitoring Data
Detected Compounds Only

| LOCATION | | | MW-1 | | | | | | | | | |
|---------------------------------------|---------|-------|-----------|----------|-----------|------------|------------|-----------|-----------|-----------|-----------|----------|
| SAMPLING DATE | | | 3/27/2018 | 6/6/2018 | 9/26/2018 | 12/19/2018 | 12/19/2018 | 3/20/2019 | 3/20/2019 | 8/23/2019 | 8/23/2019 | 4/9/2020 |
| | | | DUPLICATE | | DUPLICATE | | DUPLICATE | | DUPLICATE | | | |
| | | | Results | | | | | | | | | |
| | NY-AWQS | Units | | | | | | | | | | |
| Dissolved Metals | | | | | | | | | | | | |
| Aluminum, Dissolved | NS | ug/l | 6.81 J | 9.78 J | 46.9 | 22 J | 604 J | 324 J | 234 J | 1580 J | 310 J | NA |
| Antimony, Dissolved | 3 | ug/l | 2.78 J | 0.54 J | 0.57 J | 1.2 J | 0.76 J | 0.84 J+ | 1.32 J+ | 1.58 J | 1.44 J | NA |
| Arsenic, Dissolved | 25 | ug/l | 6.8 | 3.71 | 1.83 | 2.18 | 2.38 | 3.33 | 3.24 | 3.65 | 3.2 | NA |
| Barium, Dissolved | 1000 | ug/l | 185.2 | 53.62 | 215.8 | 149.8 | 148.1 | 19.97 | 17.12 | 30.06 J | 17.28 J | NA |
| Cadmium, Dissolved | 5 | ug/l | | 0.28 | 0.22 | 0.11 J | 0.1 J | | | | | NA |
| Calcium, Dissolved | NS | ug/l | 57600 | 25100 | 146000 | 166000 | 165000 | 12000 | 11700 | 7730 | 6720 | NA |
| Chromium, Dissolved | 50 | ug/l | | | | | 0.74 J | 0.43 J | 0.33 J | 1.81 | 0.19 J | NA |
| Cobalt, Dissolved | NS | ug/l | 1.1 | 0.41 J | 1.44 | 1.6 J | 2.02 J | 0.31 J | 0.22 J | 1.05 | 0.25 J | NA |
| Copper, Dissolved | 200 | ug/l | 1.13 | 0.99 J | 1.25 | 0.93 J | 2.21 | 1.37 | 1.21 | 3.78 J | 1.39 J | NA |
| Iron, Dissolved | 300 | ug/l | | | 83.2 | 67.9 J+ | 998 J+ | 385 J | 193 J | 1860 J+ | 293 J+ | NA |
| Lead, Dissolved | 25 | ug/l | | | | | 0.73 J | | | 1.22 | | NA |
| Magnesium, Dissolved | 35000 | ug/l | 42300 | 16200 | 103000 | 141000 | 141000 J | 7290 | 7140 | 4290 | 3550 | NA |
| Manganese, Dissolved | 300 | ug/l | 235.2 | 152.8 | 967.2 | 1041 | 1032 | 54.8 | 51.1 | 56.8 J | 29.48 J | NA |
| Mercury, Dissolved | 0.7 | ug/l | | | 0.2 UJ | | | | | | | NA |
| Nickel, Dissolved | 100 | ug/l | 2.55 | 1.21 J | 1.67 J | 2.15 | 2.59 | 0.73 J | | 3.5 | 1.9 J | NA |
| Potassium, Dissolved | NS | ug/l | 2080 | 871 | 2490 | 2300 | 2580 | 822 | 759 | 1120 J | 633 J | NA |
| Silver, Dissolved | 50 | ug/l | | | 0.22 J | | | | | | | NA |
| Sodium, Dissolved | 20000 | ug/l | 97500 | 73400 | 193000 | 217000 | 214000 | 68200 | 68100 | 68700 | 65100 | NA |
| Thallium, Dissolved | 0.5 | ug/l | | | | | | | | | 0.19 J | NA |
| Vanadium, Dissolved | NS | ug/l | 2.15 J | | | | 1.85 J | 2.13 J | 1.82 J | 6.77 | 4.42 J | NA |
| Zinc, Dissolved | 2000 | ug/l | | | 4.41 J | | | | | 8.38 J | | NA |
| Total Metals | | | | | | | | | | | | |
| Aluminum, Total | NS | ug/l | 3790 | 7650 | 2080 | 16100 J | 20800 J | 11700 | 9800 | 12000 | 11900 | NA |
| Antimony, Total | 3 | ug/l | 2.23 J | 0.56 J | | 0.92 J | 1.95 J | 0.92 J | 1.38 J | | 0.47 J | NA |
| Arsenic, Total | 25 | ug/l | 9.06 | 6.7 | 3.3 | 18.12 J | 24.9 J | 12.27 | 10.76 | 12.09 | 12.18 | NA |
| Barium, Total | 1000 | ug/l | 228.5 | 99.23 | 244.4 | 511.3 | 518.6 | 98.99 | 86.21 | 125.8 | 127.2 | NA |
| Beryllium, Total | 3 | ug/l | 0.21 J | 0.36 J | 0.15 J | 1.51 | 1.72 | 0.58 | 0.49 J | 0.83 J | 0.64 J | NA |
| Cadmium, Total | 5 | ug/l | 0.06 J | 0.57 | 0.47 | 1.14 | 1.12 | 0.13 J | 0.14 J | 0.19 J | 0.22 | NA |
| Calcium, Total | NS | ug/l | 75500 | 42400 | 167000 | 820000 | 844000 | 57200 | 54000 | 69500 | 78100 | NA |
| Chromium, Total | 50 | ug/l | 5.3 | 11.2 | 3.67 | 29.51 J | 39.21 J | 16.91 | 14.59 | 21.48 | 20.49 | NA |
| Cobalt, Total | NS | ug/l | 4.29 | 6.63 | 4 | 36.94 J | 46.53 J | 14.23 J | 10.83 J | 14.64 | 14.69 | NA |
| Copper, Total | 200 | ug/l | 9 | 20.73 | 11.33 | 201.6 | 226.6 | 45.23 J | 35.41 J | 48.39 | 52.77 | NA |
| Iron, Total | 300 | ug/l | 7290 | 14400 | 4980 J+ | 88400 | 107000 | 27700 J+ | 21500 J+ | 31600 J- | 31700 J- | NA |
| Lead, Total | 25 | ug/l | 2.96 | 6.85 | 4.4 | 53.64 | 63.19 | 17.36 | 14.23 | 17.3 | 17.89 | NA |
| Magnesium, Total | 35000 | ug/l | 46100 | 23100 | 113000 J+ | 284000 | 274000 | 21600 | 19500 | 23800 | 25500 | NA |
| Manganese, Total | 300 | ug/l | 417 | 468 | 1270 | 10080 | 10390 | 869 | 784.5 | 1053 | 1131 | NA |
| Nickel, Total | 100 | ug/l | 9.82 | 13.88 | 6 | 57.34 J | 77.78 J | 29.38 J | 23.4 J | 31.6 | 33.02 | NA |
| Potassium, Total | NS | ug/l | 3230 | 2410 | 2930 | 6540 J | 8910 J | 2590 | 2330 | 2900 | 2820 | NA |
| Selenium, Total | 10 | ug/l | | | | 8.83 | 9.54 | 3.37 J | 3.16 J | 5.71 | 5.33 | NA |
| Silver, Total | 50 | ug/l | | 0.19 J | | | 0.21 J | | | | | NA |
| Sodium, Total | 20000 | ug/l | 99500 | 73100 | 200000 | 204000 | 195000 | 73400 | 74600 | 72900 | 75900 | NA |
| Thallium, Total | 0.5 | ug/l | | | | 0.28 J | 0.48 J | 0.16 J | 0.15 J | 0.26 JR | 0.29 J | NA |
| Vanadium, Total | NS | ug/l | 9.07 | 12.62 | 5.75 | 47.17 | 55.38 | 21.65 J | 17.37 J | 27.04 | 26.61 | NA |
| Zinc, Total | 2000 | ug/l | 21.91 | 45.51 | 22.44 | 245 J | 320.4 J | 72.71 | 60.32 | 88.72 | 90.72 | NA |
| Volatile Organics by GC/MS | | | | | | | | | | | | |
| 1,4-Dioxane | | ug/l | | | 250 UR | | 61 UR | | | | | |
| Acetone | 50 | ug/l | | | | | | | | | | 17 |
| Bromomethane | 5 | ug/l | | | 2.5 UJ | | | | | | | |
| Methyl tert butyl ether | 10 | ug/l | 2300 | 660 | 3700 | 2800 | 2300 | 390 | 390 | 140 | 150 | 34 |
| Tetrachloroethene | 5 | ug/l | | 0.59 J | | | | | | | | |
| Semivolatile Organics by GC/MS | | | | | | | | | | | | |
| 3,3'-Dichlorobenzidine | 5 | ug/l | | | | 0.5 UJ | 0.5 UJ | | | | | |
| 4-Chloroaniline | 5 | ug/l | | | 5 UJ | 1.1 UJ | 1.1 UJ | | | | | |
| 4-Nitroaniline | 5 | ug/l | | | | 0.5 UJ | 0.5 UJ | | | | | |
| Acetophenone | | ug/l | | | | | | | | | | |
| Bis(2-ethylhexyl)phthalate | 5 | ug/l | 3.7 | | | 52 J+ | | | | | | |
| Di-n-butylphthalate | 50 | ug/l | | | | 0.84 J | | | | | | |
| 2-Chloronaphthalene | 10 | ug/l | | | | | | | | | | |
| 2-Methylnaphthalene | NS | ug/l | | | | | | 0.02 J | | | | |
| Anthracene | 50 | ug/l | 0.1 | | | | | | | | | |
| Fluoranthene | 50 | ug/l | 0.1 | | | | 0.03 J | | | | | |
| Fluorene | 50 | ug/l | 0.1 | | | | | | | | | |
| Naphthalene | 10 | ug/l | 0.1 | | | | | | | | 0.42 BR | |
| Phenanthrene | 50 | ug/l | 0.1 | | | | | 0.04 JU | 0.05 JU | | | |
| Pyrene | 5 | ug/l | | | | | 0.02 J | | | | | |

Notes:
 NY-AWQS: New York TOGS 111 Ambient Water Quality Standards criteria reflects all addendum to criteria through June 2004.
 Samples were analyzed by Alpha Analytical Laboratory
 Blank cells indicate parameters that were not detected
 Shaded values indicate a value greater than NY TOGS 1.1.1 Ambient Water Quality Standards Limits
 R - indicates a rejected, unusable value dictated by the DUSR report
 J+ - indicates an estimated, biased high value dictated by the DUSR report
 J - indicates an estimated value
 J- - indicates an estimated, biased low value dictated by the DUSR report
 UJ - indicates an estimated value dictated by the DUSR report
 NS - no standard
 NA - Not Analyzed

APPENDIX A
MONITORING WELL SAMPLING LOG

APPENDIX B
LABORATORY ANALYTICAL DATA



ANALYTICAL REPORT

| | |
|-----------------|---|
| Lab Number: | L2015218 |
| Client: | CHA Companies 3 Winners Circle Albany, NY 12205 |
| ATTN: | C. Hurlburt |
| Phone: | (518) 453-4500 |
| Project Name: | Not Specified |
| Project Number: | 31647 |
| Report Date: | 04/16/20 |

The original project report/data package is held by Alpha Analytical. This report/data package is paginated and should be reproduced only in its entirety. Alpha Analytical holds no responsibility for results and/or data that are not consistent with the original.

Certifications & Approvals: MA (M-MA086), NH NELAP (2064), CT (PH-0574), IL (200077), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #P330-17-00196).

Eight Walkup Drive, Westborough, MA 01581-1019
508-898-9220 (Fax) 508-898-9193 800-624-9220 - www.alphalab.com



Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

| Alpha Sample ID | Client ID | Matrix | Sample Location | Collection Date/Time | Receive Date |
|----------------------------|------------------|---------------|----------------------------|---------------------------------|---------------------|
| L2015218-01 | GW-MW1-040920 | WATER | Not Specified | 04/09/20 11:45 | 04/09/20 |
| L2015218-02 | TRIP BLANK | WATER | Not Specified | 04/09/20 00:00 | 04/09/20 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Case Narrative

The samples were received in accordance with the Chain of Custody and no significant deviations were encountered during the preparation or analysis unless otherwise noted. Sample Receipt, Container Information, and the Chain of Custody are located at the back of the report.

Results contained within this report relate only to the samples submitted under this Alpha Lab Number and meet NELAP requirements for all NELAP accredited parameters unless otherwise noted in the following narrative. The data presented in this report is organized by parameter (i.e. VOC, SVOC, etc.). Sample specific Quality Control data (i.e. Surrogate Spike Recovery) is reported at the end of the target analyte list for each individual sample, followed by the Laboratory Batch Quality Control at the end of each parameter. Tentatively Identified Compounds (TICs), if requested, are reported for compounds identified to be present and are not part of the method/program Target Compound List, even if only a subset of the TCL are being reported. If a sample was re-analyzed or re-extracted due to a required quality control corrective action and if both sets of data are reported, the Laboratory ID of the re-analysis or re-extraction is designated with an "R" or "RE", respectively.

When multiple Batch Quality Control elements are reported (e.g. more than one LCS), the associated samples for each element are noted in the grey shaded header line of each data table. Any Laboratory Batch, Sample Specific % recovery or RPD value that is outside the listed Acceptance Criteria is bolded in the report. In reference to questions H (CAM) or 4 (RCP) when "NO" is checked, the performance criteria for CAM and RCP methods allow for some quality control failures to occur and still be within method compliance. In these instances, the specific failure is not narrated but noted in the associated QC Outlier Summary Report, located directly after the Case Narrative. QC information is also incorporated in the Data Usability Assessment table (Format 11) of our Data Merger tool, where it can be reviewed in conjunction with the sample result, associated regulatory criteria and any associated data usability implications.

Soil/sediments, solids and tissues are reported on a dry weight basis unless otherwise noted. Definitions of all data qualifiers and acronyms used in this report are provided in the Glossary located at the back of the report.

HOLD POLICY - For samples submitted on hold, Alpha's policy is to hold samples (with the exception of Air canisters) free of charge for 21 calendar days from the date the project is completed. After 21 calendar days, we will dispose of all samples submitted including those put on hold unless you have contacted your Alpha Project Manager and made arrangements for Alpha to continue to hold the samples. Air canisters will be disposed after 3 business days from the date the project is completed.

Please contact Project Management at 800-624-9220 with any questions.

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Case Narrative (continued)

Report Submission

All non-detect (ND) or estimated concentrations (J-qualified) have been quantitated to the limit noted in the MDL column.

Sample Receipt

L2015218-02: A sample identified as "TRIP BLANK" was received, but not listed on the Chain of Custody. This sample was not analyzed.

I, the undersigned, attest under the pains and penalties of perjury that, to the best of my knowledge and belief and based upon my personal inquiry of those responsible for providing the information contained in this analytical report, such information is accurate and complete. This certificate of analysis is not complete unless this page accompanies any and all pages of this report.

Authorized Signature:



Michelle M. Morris

Title: Technical Director/Representative

Date: 04/16/20

ORGANICS

VOLATILES

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

SAMPLE RESULTS

Lab ID: L2015218-01
 Client ID: GW-MW1-040920
 Sample Location: Not Specified

Date Collected: 04/09/20 11:45
 Date Received: 04/09/20
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 04/14/20 01:13
 Analyst: NLK

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|------|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab | | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 | 1 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 | 1 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 | 1 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 | 1 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 | 1 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 | 1 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 | 1 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 | 1 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 | 1 |
| Benzene | ND | | ug/l | 0.50 | 0.16 | 1 |
| Toluene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Chloromethane | 1.6 | J | ug/l | 2.5 | 0.70 | 1 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 | 1 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 | 1 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Trichloroethene | ND | | ug/l | 0.50 | 0.18 | 1 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

SAMPLE RESULTS

Lab ID: L2015218-01
 Client ID: GW-MW1-040920
 Sample Location: Not Specified

Date Collected: 04/09/20 11:45
 Date Received: 04/09/20
 Field Prep: Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------|-----------|-------|-----|------|-----------------|
| Volatile Organics by GC/MS - Westborough Lab | | | | | | |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl tert butyl ether | 34 | | ug/l | 2.5 | 0.70 | 1 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 | 1 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Styrene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 | 1 |
| Acetone | 17 | | ug/l | 5.0 | 1.5 | 1 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 | 1 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 | 1 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 | 1 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 | 1 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl Acetate | ND | | ug/l | 2.0 | 0.23 | 1 |
| Cyclohexane | ND | | ug/l | 10 | 0.27 | 1 |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. | 1 |
| Freon-113 | ND | | ug/l | 2.5 | 0.70 | 1 |
| Methyl cyclohexane | ND | | ug/l | 10 | 0.40 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|-----------------------|------------|-----------|---------------------|
| 1,2-Dichloroethane-d4 | 116 | | 70-130 |
| Toluene-d8 | 99 | | 70-130 |
| 4-Bromofluorobenzene | 90 | | 70-130 |
| Dibromofluoromethane | 107 | | 70-130 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 04/13/20 17:43
Analyst: AD

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|------|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1360924-5 | | | | | |
| Methylene chloride | ND | | ug/l | 2.5 | 0.70 |
| 1,1-Dichloroethane | ND | | ug/l | 2.5 | 0.70 |
| Chloroform | ND | | ug/l | 2.5 | 0.70 |
| Carbon tetrachloride | ND | | ug/l | 0.50 | 0.13 |
| 1,2-Dichloropropane | ND | | ug/l | 1.0 | 0.14 |
| Dibromochloromethane | ND | | ug/l | 0.50 | 0.15 |
| 1,1,2-Trichloroethane | ND | | ug/l | 1.5 | 0.50 |
| Tetrachloroethene | ND | | ug/l | 0.50 | 0.18 |
| Chlorobenzene | ND | | ug/l | 2.5 | 0.70 |
| Trichlorofluoromethane | ND | | ug/l | 2.5 | 0.70 |
| 1,2-Dichloroethane | ND | | ug/l | 0.50 | 0.13 |
| 1,1,1-Trichloroethane | ND | | ug/l | 2.5 | 0.70 |
| Bromodichloromethane | ND | | ug/l | 0.50 | 0.19 |
| trans-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.16 |
| cis-1,3-Dichloropropene | ND | | ug/l | 0.50 | 0.14 |
| Bromoform | ND | | ug/l | 2.0 | 0.65 |
| 1,1,2,2-Tetrachloroethane | ND | | ug/l | 0.50 | 0.17 |
| Benzene | ND | | ug/l | 0.50 | 0.16 |
| Toluene | ND | | ug/l | 2.5 | 0.70 |
| Ethylbenzene | ND | | ug/l | 2.5 | 0.70 |
| Chloromethane | ND | | ug/l | 2.5 | 0.70 |
| Bromomethane | ND | | ug/l | 2.5 | 0.70 |
| Vinyl chloride | ND | | ug/l | 1.0 | 0.07 |
| Chloroethane | ND | | ug/l | 2.5 | 0.70 |
| 1,1-Dichloroethene | ND | | ug/l | 0.50 | 0.17 |
| trans-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 |
| Trichloroethene | ND | | ug/l | 0.50 | 0.18 |
| 1,2-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 |
| 1,3-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 04/13/20 17:43
Analyst: AD

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|-----|------|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1360924-5 | | | | | |
| 1,4-Dichlorobenzene | ND | | ug/l | 2.5 | 0.70 |
| Methyl tert butyl ether | ND | | ug/l | 2.5 | 0.70 |
| p/m-Xylene | ND | | ug/l | 2.5 | 0.70 |
| o-Xylene | ND | | ug/l | 2.5 | 0.70 |
| cis-1,2-Dichloroethene | ND | | ug/l | 2.5 | 0.70 |
| Styrene | ND | | ug/l | 2.5 | 0.70 |
| Dichlorodifluoromethane | ND | | ug/l | 5.0 | 1.0 |
| Acetone | ND | | ug/l | 5.0 | 1.5 |
| Carbon disulfide | ND | | ug/l | 5.0 | 1.0 |
| 2-Butanone | ND | | ug/l | 5.0 | 1.9 |
| 4-Methyl-2-pentanone | ND | | ug/l | 5.0 | 1.0 |
| 2-Hexanone | ND | | ug/l | 5.0 | 1.0 |
| Bromochloromethane | ND | | ug/l | 2.5 | 0.70 |
| 1,2-Dibromoethane | ND | | ug/l | 2.0 | 0.65 |
| 1,2-Dibromo-3-chloropropane | ND | | ug/l | 2.5 | 0.70 |
| Isopropylbenzene | ND | | ug/l | 2.5 | 0.70 |
| 1,2,3-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 |
| 1,2,4-Trichlorobenzene | ND | | ug/l | 2.5 | 0.70 |
| Methyl Acetate | ND | | ug/l | 2.0 | 0.23 |
| Cyclohexane | ND | | ug/l | 10 | 0.27 |
| 1,4-Dioxane | ND | | ug/l | 250 | 61. |
| Freon-113 | ND | | ug/l | 2.5 | 0.70 |
| Methyl cyclohexane | ND | | ug/l | 10 | 0.40 |

Project Name: Not Specified**Lab Number:** L2015218**Project Number:** 31647**Report Date:** 04/16/20

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
 Analytical Date: 04/13/20 17:43
 Analyst: AD

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|----|-----|
| Volatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1360924-5 | | | | | |

| Surrogate | %Recovery | Qualifier | Acceptance Criteria |
|-----------------------|-----------|-----------|------------------------|
| 1,2-Dichloroethane-d4 | 111 | | 70-130 |
| Toluene-d8 | 101 | | 70-130 |
| 4-Bromofluorobenzene | 95 | | 70-130 |
| Dibromofluoromethane | 97 | | 70-130 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1360924-3 WG1360924-4 | | | | | | | | |
| Methylene chloride | 120 | | 110 | | 70-130 | 9 | | 20 |
| 1,1-Dichloroethane | 130 | | 120 | | 70-130 | 8 | | 20 |
| Chloroform | 110 | | 110 | | 70-130 | 0 | | 20 |
| Carbon tetrachloride | 99 | | 95 | | 63-132 | 4 | | 20 |
| 1,2-Dichloropropane | 120 | | 110 | | 70-130 | 9 | | 20 |
| Dibromochloromethane | 100 | | 100 | | 63-130 | 0 | | 20 |
| 1,1,2-Trichloroethane | 110 | | 110 | | 70-130 | 0 | | 20 |
| Tetrachloroethene | 97 | | 94 | | 70-130 | 3 | | 20 |
| Chlorobenzene | 100 | | 100 | | 75-130 | 0 | | 20 |
| Trichlorofluoromethane | 130 | | 120 | | 62-150 | 8 | | 20 |
| 1,2-Dichloroethane | 120 | | 120 | | 70-130 | 0 | | 20 |
| 1,1,1-Trichloroethane | 96 | | 100 | | 67-130 | 4 | | 20 |
| Bromodichloromethane | 110 | | 110 | | 67-130 | 0 | | 20 |
| trans-1,3-Dichloropropene | 93 | | 93 | | 70-130 | 0 | | 20 |
| cis-1,3-Dichloropropene | 100 | | 100 | | 70-130 | 0 | | 20 |
| Bromoform | 93 | | 97 | | 54-136 | 4 | | 20 |
| 1,1,2,2-Tetrachloroethane | 110 | | 110 | | 67-130 | 0 | | 20 |
| Benzene | 110 | | 110 | | 70-130 | 0 | | 20 |
| Toluene | 100 | | 100 | | 70-130 | 0 | | 20 |
| Ethylbenzene | 110 | | 100 | | 70-130 | 10 | | 20 |
| Chloromethane | 140 | Q | 130 | | 64-130 | 7 | | 20 |
| Bromomethane | 95 | | 99 | | 39-139 | 4 | | 20 |
| Vinyl chloride | 130 | | 120 | | 55-140 | 8 | | 20 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS | Qual | LCS | Qual | %Recovery | RPD | Qual | RPD |
|--|-----------|------|-----------|------|-----------|-----|------|--------|
| | %Recovery | | %Recovery | | Limits | | | Limits |
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1360924-3 WG1360924-4 | | | | | | | | |
| Chloroethane | 150 | Q | 140 | Q | 55-138 | 7 | | 20 |
| 1,1-Dichloroethene | 110 | | 100 | | 61-145 | 10 | | 20 |
| trans-1,2-Dichloroethene | 110 | | 100 | | 70-130 | 10 | | 20 |
| Trichloroethene | 110 | | 110 | | 70-130 | 0 | | 20 |
| 1,2-Dichlorobenzene | 100 | | 98 | | 70-130 | 2 | | 20 |
| 1,3-Dichlorobenzene | 100 | | 99 | | 70-130 | 1 | | 20 |
| 1,4-Dichlorobenzene | 100 | | 95 | | 70-130 | 5 | | 20 |
| Methyl tert butyl ether | 97 | | 110 | | 63-130 | 13 | | 20 |
| p/m-Xylene | 110 | | 105 | | 70-130 | 5 | | 20 |
| o-Xylene | 105 | | 105 | | 70-130 | 0 | | 20 |
| cis-1,2-Dichloroethene | 110 | | 110 | | 70-130 | 0 | | 20 |
| Styrene | 105 | | 100 | | 70-130 | 5 | | 20 |
| Dichlorodifluoromethane | 96 | | 93 | | 36-147 | 3 | | 20 |
| Acetone | 130 | | 130 | | 58-148 | 0 | | 20 |
| Carbon disulfide | 120 | | 110 | | 51-130 | 9 | | 20 |
| 2-Butanone | 130 | | 130 | | 63-138 | 0 | | 20 |
| 4-Methyl-2-pentanone | 110 | | 120 | | 59-130 | 9 | | 20 |
| 2-Hexanone | 110 | | 120 | | 57-130 | 9 | | 20 |
| Bromochloromethane | 110 | | 110 | | 70-130 | 0 | | 20 |
| 1,2-Dibromoethane | 100 | | 100 | | 70-130 | 0 | | 20 |
| 1,2-Dibromo-3-chloropropane | 93 | | 92 | | 41-144 | 1 | | 20 |
| Isopropylbenzene | 100 | | 96 | | 70-130 | 4 | | 20 |
| 1,2,3-Trichlorobenzene | 92 | | 92 | | 70-130 | 0 | | 20 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Project Number: 31647

Lab Number: L2015218

Report Date: 04/16/20

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1360924-3 WG1360924-4 | | | | | | | | |
| 1,2,4-Trichlorobenzene | 89 | | 89 | | 70-130 | 0 | | 20 |
| Methyl Acetate | 120 | | 130 | | 70-130 | 8 | | 20 |
| Cyclohexane | 110 | | 110 | | 70-130 | 0 | | 20 |
| 1,4-Dioxane | 154 | | 158 | | 56-162 | 3 | | 20 |
| Freon-113 | 110 | | 100 | | 70-130 | 10 | | 20 |
| Methyl cyclohexane | 110 | | 100 | | 70-130 | 10 | | 20 |

| Surrogate | LCS %Recovery | Qual | LCSD %Recovery | Qual | Acceptance Criteria |
|-----------------------|------------------|------|-------------------|------|------------------------|
| 1,2-Dichloroethane-d4 | 111 | | 111 | | 70-130 |
| Toluene-d8 | 101 | | 99 | | 70-130 |
| 4-Bromofluorobenzene | 94 | | 91 | | 70-130 |
| Dibromofluoromethane | 102 | | 102 | | 70-130 |

SEMIVOLATILES

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

SAMPLE RESULTS

Lab ID: L2015218-01
 Client ID: GW-MW1-040920
 Sample Location: Not Specified

Date Collected: 04/09/20 11:45
 Date Received: 04/09/20
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8270D
 Analytical Date: 04/16/20 00:22
 Analyst: WR

Extraction Method: EPA 3510C
 Extraction Date: 04/15/20 15:24

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab | | | | | | |
| Bis(2-chloroethyl)ether | ND | | ug/l | 2.0 | 0.50 | 1 |
| 3,3'-Dichlorobenzidine | ND | | ug/l | 5.0 | 1.6 | 1 |
| 2,4-Dinitrotoluene | ND | | ug/l | 5.0 | 1.2 | 1 |
| 2,6-Dinitrotoluene | ND | | ug/l | 5.0 | 0.93 | 1 |
| 4-Chlorophenyl phenyl ether | ND | | ug/l | 2.0 | 0.49 | 1 |
| 4-Bromophenyl phenyl ether | ND | | ug/l | 2.0 | 0.38 | 1 |
| Bis(2-chloroisopropyl)ether | ND | | ug/l | 2.0 | 0.53 | 1 |
| Bis(2-chloroethoxy)methane | ND | | ug/l | 5.0 | 0.50 | 1 |
| Hexachlorocyclopentadiene | ND | | ug/l | 20 | 0.69 | 1 |
| Isophorone | ND | | ug/l | 5.0 | 1.2 | 1 |
| Nitrobenzene | ND | | ug/l | 2.0 | 0.77 | 1 |
| NDPA/DPA | ND | | ug/l | 2.0 | 0.42 | 1 |
| n-Nitrosodi-n-propylamine | ND | | ug/l | 5.0 | 0.64 | 1 |
| Bis(2-ethylhexyl)phthalate | ND | | ug/l | 3.0 | 1.5 | 1 |
| Butyl benzyl phthalate | ND | | ug/l | 5.0 | 1.2 | 1 |
| Di-n-butylphthalate | ND | | ug/l | 5.0 | 0.39 | 1 |
| Di-n-octylphthalate | ND | | ug/l | 5.0 | 1.3 | 1 |
| Diethyl phthalate | ND | | ug/l | 5.0 | 0.38 | 1 |
| Dimethyl phthalate | ND | | ug/l | 5.0 | 1.8 | 1 |
| Biphenyl | ND | | ug/l | 2.0 | 0.46 | 1 |
| 4-Chloroaniline | ND | | ug/l | 5.0 | 1.1 | 1 |
| 2-Nitroaniline | ND | | ug/l | 5.0 | 0.50 | 1 |
| 3-Nitroaniline | ND | | ug/l | 5.0 | 0.81 | 1 |
| 4-Nitroaniline | ND | | ug/l | 5.0 | 0.80 | 1 |
| Dibenzofuran | ND | | ug/l | 2.0 | 0.50 | 1 |
| 1,2,4,5-Tetrachlorobenzene | ND | | ug/l | 10 | 0.44 | 1 |
| Acetophenone | ND | | ug/l | 5.0 | 0.53 | 1 |
| 2,4,6-Trichlorophenol | ND | | ug/l | 5.0 | 0.61 | 1 |

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

SAMPLE RESULTS

Lab ID: L2015218-01
 Client ID: GW-MW1-040920
 Sample Location: Not Specified

Date Collected: 04/09/20 11:45
 Date Received: 04/09/20
 Field Prep: Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|---|--------|-----------|-------|-----|------|-----------------|
| Semivolatile Organics by GC/MS - Westborough Lab | | | | | | |
| p-Chloro-m-cresol | ND | | ug/l | 2.0 | 0.35 | 1 |
| 2-Chlorophenol | ND | | ug/l | 2.0 | 0.48 | 1 |
| 2,4-Dichlorophenol | ND | | ug/l | 5.0 | 0.41 | 1 |
| 2,4-Dimethylphenol | ND | | ug/l | 5.0 | 1.8 | 1 |
| 2-Nitrophenol | ND | | ug/l | 10 | 0.85 | 1 |
| 4-Nitrophenol | ND | | ug/l | 10 | 0.67 | 1 |
| 2,4-Dinitrophenol | ND | | ug/l | 20 | 6.6 | 1 |
| 4,6-Dinitro-o-cresol | ND | | ug/l | 10 | 1.8 | 1 |
| Phenol | ND | | ug/l | 5.0 | 0.57 | 1 |
| 2-Methylphenol | ND | | ug/l | 5.0 | 0.49 | 1 |
| 3-Methylphenol/4-Methylphenol | ND | | ug/l | 5.0 | 0.48 | 1 |
| 2,4,5-Trichlorophenol | ND | | ug/l | 5.0 | 0.77 | 1 |
| Carbazole | ND | | ug/l | 2.0 | 0.49 | 1 |
| Atrazine | ND | | ug/l | 10 | 0.76 | 1 |
| Benzaldehyde | ND | | ug/l | 5.0 | 0.53 | 1 |
| Caprolactam | ND | | ug/l | 10 | 3.3 | 1 |
| 2,3,4,6-Tetrachlorophenol | ND | | ug/l | 5.0 | 0.84 | 1 |

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol | 54 | | 21-120 |
| Phenol-d6 | 46 | | 10-120 |
| Nitrobenzene-d5 | 65 | | 23-120 |
| 2-Fluorobiphenyl | 61 | | 15-120 |
| 2,4,6-Tribromophenol | 41 | | 10-120 |
| 4-Terphenyl-d14 | 63 | | 41-149 |

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

SAMPLE RESULTS

Lab ID: L2015218-01
 Client ID: GW-MW1-040920
 Sample Location: Not Specified

Date Collected: 04/09/20 11:45
 Date Received: 04/09/20
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8270D-SIM
 Analytical Date: 04/15/20 12:19
 Analyst: CB

Extraction Method: EPA 3510C
 Extraction Date: 04/13/20 23:55

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|--|--------|-----------|-------|------|------|-----------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab | | | | | | |
| Acenaphthene | ND | | ug/l | 0.10 | 0.01 | 1 |
| 2-Chloronaphthalene | ND | | ug/l | 0.20 | 0.02 | 1 |
| Fluoranthene | ND | | ug/l | 0.10 | 0.02 | 1 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 0.05 | 1 |
| Naphthalene | ND | | ug/l | 0.10 | 0.05 | 1 |
| Benzo(a)anthracene | ND | | ug/l | 0.10 | 0.02 | 1 |
| Benzo(a)pyrene | ND | | ug/l | 0.10 | 0.02 | 1 |
| Benzo(b)fluoranthene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Benzo(k)fluoranthene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Chrysene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Acenaphthylene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Anthracene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Benzo(ghi)perylene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Fluorene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Phenanthrene | ND | | ug/l | 0.10 | 0.02 | 1 |
| Dibenzo(a,h)anthracene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Indeno(1,2,3-cd)pyrene | ND | | ug/l | 0.10 | 0.01 | 1 |
| Pyrene | ND | | ug/l | 0.10 | 0.02 | 1 |
| 2-Methylnaphthalene | ND | | ug/l | 0.10 | 0.02 | 1 |
| Pentachlorophenol | ND | | ug/l | 0.80 | 0.01 | 1 |
| Hexachlorobenzene | ND | | ug/l | 0.80 | 0.01 | 1 |
| Hexachloroethane | ND | | ug/l | 0.80 | 0.06 | 1 |

Project Name: Not Specified**Lab Number:** L2015218**Project Number:** 31647**Report Date:** 04/16/20**SAMPLE RESULTS**

Lab ID: L2015218-01

Date Collected: 04/09/20 11:45

Client ID: GW-MW1-040920

Date Received: 04/09/20

Sample Location: Not Specified

Field Prep: Not Specified

Sample Depth:

| Parameter | Result | Qualifier | Units | RL | MDL | Dilution Factor |
|-----------|--------|-----------|-------|----|-----|-----------------|
|-----------|--------|-----------|-------|----|-----|-----------------|

Semivolatile Organics by GC/MS-SIM - Westborough Lab

| Surrogate | % Recovery | Qualifier | Acceptance Criteria |
|----------------------|------------|-----------|---------------------|
| 2-Fluorophenol | 53 | | 21-120 |
| Phenol-d6 | 48 | | 10-120 |
| Nitrobenzene-d5 | 66 | | 23-120 |
| 2-Fluorobiphenyl | 62 | | 15-120 |
| 2,4,6-Tribromophenol | 50 | | 10-120 |
| 4-Terphenyl-d14 | 63 | | 41-149 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D-SIM
Analytical Date: 04/15/20 10:55
Analyst: JJW

Extraction Method: EPA 3510C
Extraction Date: 04/13/20 23:55

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|------|------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01 Batch: WG1360851-1 | | | | | |
| Acenaphthene | ND | | ug/l | 0.10 | 0.01 |
| 2-Chloronaphthalene | ND | | ug/l | 0.20 | 0.02 |
| Fluoranthene | ND | | ug/l | 0.10 | 0.02 |
| Hexachlorobutadiene | ND | | ug/l | 0.50 | 0.05 |
| Naphthalene | ND | | ug/l | 0.10 | 0.05 |
| Benzo(a)anthracene | ND | | ug/l | 0.10 | 0.02 |
| Benzo(a)pyrene | ND | | ug/l | 0.10 | 0.02 |
| Benzo(b)fluoranthene | ND | | ug/l | 0.10 | 0.01 |
| Benzo(k)fluoranthene | ND | | ug/l | 0.10 | 0.01 |
| Chrysene | ND | | ug/l | 0.10 | 0.01 |
| Acenaphthylene | ND | | ug/l | 0.10 | 0.01 |
| Anthracene | ND | | ug/l | 0.10 | 0.01 |
| Benzo(ghi)perylene | ND | | ug/l | 0.10 | 0.01 |
| Fluorene | ND | | ug/l | 0.10 | 0.01 |
| Phenanthrene | ND | | ug/l | 0.10 | 0.02 |
| Dibenzo(a,h)anthracene | ND | | ug/l | 0.10 | 0.01 |
| Indeno(1,2,3-cd)pyrene | ND | | ug/l | 0.10 | 0.01 |
| Pyrene | ND | | ug/l | 0.10 | 0.02 |
| 2-Methylnaphthalene | ND | | ug/l | 0.10 | 0.02 |
| Pentachlorophenol | ND | | ug/l | 0.80 | 0.01 |
| Hexachlorobenzene | ND | | ug/l | 0.80 | 0.01 |
| Hexachloroethane | ND | | ug/l | 0.80 | 0.06 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D-SIM
Analytical Date: 04/15/20 10:55
Analyst: JJW

Extraction Method: EPA 3510C
Extraction Date: 04/13/20 23:55

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|----|-----|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01 Batch: WG1360851-1 | | | | | |

| Surrogate | %Recovery | Qualifier | Acceptance Criteria |
|----------------------|-----------|-----------|---------------------|
| 2-Fluorophenol | 42 | | 21-120 |
| Phenol-d6 | 41 | | 10-120 |
| Nitrobenzene-d5 | 58 | | 23-120 |
| 2-Fluorobiphenyl | 57 | | 15-120 |
| 2,4,6-Tribromophenol | 33 | | 10-120 |
| 4-Terphenyl-d14 | 65 | | 41-149 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D
Analytical Date: 04/15/20 23:11
Analyst: WR

Extraction Method: EPA 3510C
Extraction Date: 04/15/20 15:24

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1361431-1 | | | | | |
| Bis(2-chloroethyl)ether | ND | | ug/l | 2.0 | 0.50 |
| 3,3'-Dichlorobenzidine | ND | | ug/l | 5.0 | 1.6 |
| 2,4-Dinitrotoluene | ND | | ug/l | 5.0 | 1.2 |
| 2,6-Dinitrotoluene | ND | | ug/l | 5.0 | 0.93 |
| 4-Chlorophenyl phenyl ether | ND | | ug/l | 2.0 | 0.49 |
| 4-Bromophenyl phenyl ether | ND | | ug/l | 2.0 | 0.38 |
| Bis(2-chloroisopropyl)ether | ND | | ug/l | 2.0 | 0.53 |
| Bis(2-chloroethoxy)methane | ND | | ug/l | 5.0 | 0.50 |
| Hexachlorocyclopentadiene | ND | | ug/l | 20 | 0.69 |
| Isophorone | ND | | ug/l | 5.0 | 1.2 |
| Nitrobenzene | ND | | ug/l | 2.0 | 0.77 |
| NDPA/DPA | ND | | ug/l | 2.0 | 0.42 |
| n-Nitrosodi-n-propylamine | ND | | ug/l | 5.0 | 0.64 |
| Bis(2-ethylhexyl)phthalate | ND | | ug/l | 3.0 | 1.5 |
| Butyl benzyl phthalate | ND | | ug/l | 5.0 | 1.2 |
| Di-n-butylphthalate | ND | | ug/l | 5.0 | 0.39 |
| Di-n-octylphthalate | ND | | ug/l | 5.0 | 1.3 |
| Diethyl phthalate | ND | | ug/l | 5.0 | 0.38 |
| Dimethyl phthalate | ND | | ug/l | 5.0 | 1.8 |
| Biphenyl | ND | | ug/l | 2.0 | 0.46 |
| 4-Chloroaniline | ND | | ug/l | 5.0 | 1.1 |
| 2-Nitroaniline | ND | | ug/l | 5.0 | 0.50 |
| 3-Nitroaniline | ND | | ug/l | 5.0 | 0.81 |
| 4-Nitroaniline | ND | | ug/l | 5.0 | 0.80 |
| Dibenzofuran | ND | | ug/l | 2.0 | 0.50 |
| 1,2,4,5-Tetrachlorobenzene | ND | | ug/l | 10 | 0.44 |
| Acetophenone | ND | | ug/l | 5.0 | 0.53 |
| 2,4,6-Trichlorophenol | ND | | ug/l | 5.0 | 0.61 |
| p-Chloro-m-cresol | ND | | ug/l | 2.0 | 0.35 |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8270D
Analytical Date: 04/15/20 23:11
Analyst: WR

Extraction Method: EPA 3510C
Extraction Date: 04/15/20 15:24

| Parameter | Result | Qualifier | Units | RL | MDL |
|---|--------|-----------|-------|-----|------|
| Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1361431-1 | | | | | |
| 2-Chlorophenol | ND | | ug/l | 2.0 | 0.48 |
| 2,4-Dichlorophenol | ND | | ug/l | 5.0 | 0.41 |
| 2,4-Dimethylphenol | ND | | ug/l | 5.0 | 1.8 |
| 2-Nitrophenol | ND | | ug/l | 10 | 0.85 |
| 4-Nitrophenol | ND | | ug/l | 10 | 0.67 |
| 2,4-Dinitrophenol | ND | | ug/l | 20 | 6.6 |
| 4,6-Dinitro-o-cresol | ND | | ug/l | 10 | 1.8 |
| Phenol | ND | | ug/l | 5.0 | 0.57 |
| 2-Methylphenol | ND | | ug/l | 5.0 | 0.49 |
| 3-Methylphenol/4-Methylphenol | ND | | ug/l | 5.0 | 0.48 |
| 2,4,5-Trichlorophenol | ND | | ug/l | 5.0 | 0.77 |
| Carbazole | ND | | ug/l | 2.0 | 0.49 |
| Atrazine | ND | | ug/l | 10 | 0.76 |
| Benzaldehyde | ND | | ug/l | 5.0 | 0.53 |
| Caprolactam | ND | | ug/l | 10 | 3.3 |
| 2,3,4,6-Tetrachlorophenol | ND | | ug/l | 5.0 | 0.84 |

| Surrogate | %Recovery | Qualifier | Acceptance Criteria |
|----------------------|-----------|-----------|------------------------|
| 2-Fluorophenol | 53 | | 21-120 |
| Phenol-d6 | 42 | | 10-120 |
| Nitrobenzene-d5 | 61 | | 23-120 |
| 2-Fluorobiphenyl | 58 | | 15-120 |
| 2,4,6-Tribromophenol | 41 | | 10-120 |
| 4-Terphenyl-d14 | 59 | | 41-149 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01 Batch: WG1360851-2 WG1360851-3 | | | | | | | | |
| Acenaphthene | 53 | | 51 | | 40-140 | 4 | | 40 |
| 2-Chloronaphthalene | 52 | | 49 | | 40-140 | 6 | | 40 |
| Fluoranthene | 51 | | 54 | | 40-140 | 6 | | 40 |
| Hexachlorobutadiene | 47 | | 38 | Q | 40-140 | 21 | | 40 |
| Naphthalene | 50 | | 43 | | 40-140 | 15 | | 40 |
| Benzo(a)anthracene | 53 | | 56 | | 40-140 | 6 | | 40 |
| Benzo(a)pyrene | 56 | | 59 | | 40-140 | 5 | | 40 |
| Benzo(b)fluoranthene | 54 | | 57 | | 40-140 | 5 | | 40 |
| Benzo(k)fluoranthene | 52 | | 56 | | 40-140 | 7 | | 40 |
| Chrysene | 52 | | 56 | | 40-140 | 7 | | 40 |
| Acenaphthylene | 51 | | 49 | | 40-140 | 4 | | 40 |
| Anthracene | 53 | | 54 | | 40-140 | 2 | | 40 |
| Benzo(ghi)perylene | 57 | | 59 | | 40-140 | 3 | | 40 |
| Fluorene | 52 | | 52 | | 40-140 | 0 | | 40 |
| Phenanthrene | 53 | | 54 | | 40-140 | 2 | | 40 |
| Dibenzo(a,h)anthracene | 61 | | 62 | | 40-140 | 2 | | 40 |
| Indeno(1,2,3-cd)pyrene | 60 | | 62 | | 40-140 | 3 | | 40 |
| Pyrene | 51 | | 54 | | 40-140 | 6 | | 40 |
| 2-Methylnaphthalene | 54 | | 48 | | 40-140 | 12 | | 40 |
| Pentachlorophenol | 39 | Q | 42 | | 40-140 | 7 | | 40 |
| Hexachlorobenzene | 52 | | 52 | | 40-140 | 0 | | 40 |
| Hexachloroethane | 43 | | 34 | Q | 40-140 | 23 | | 40 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|-----------|------------------|------|-------------------|------|---------------------|-----|------|---------------|
|-----------|------------------|------|-------------------|------|---------------------|-----|------|---------------|

Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01 Batch: WG1360851-2 WG1360851-3

| Surrogate | LCS %Recovery | Qual | LCSD %Recovery | Qual | Acceptance Criteria |
|----------------------|------------------|------|-------------------|------|------------------------|
| 2-Fluorophenol | 49 | | 40 | | 21-120 |
| Phenol-d6 | 47 | | 39 | | 10-120 |
| Nitrobenzene-d5 | 57 | | 48 | | 23-120 |
| 2-Fluorobiphenyl | 52 | | 48 | | 15-120 |
| 2,4,6-Tribromophenol | 54 | | 49 | | 10-120 |
| 4-Terphenyl-d14 | 55 | | 58 | | 41-149 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS | | LCSD | | %Recovery Limits | RPD | Qual | RPD Limits |
|--|-----------|------|-----------|------|------------------|-----|------|------------|
| | %Recovery | Qual | %Recovery | Qual | | | | |
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1361431-2 WG1361431-3 | | | | | | | | |
| Bis(2-chloroethyl)ether | 64 | | 74 | | 40-140 | 14 | | 30 |
| 3,3'-Dichlorobenzidine | 37 | Q | 33 | Q | 40-140 | 11 | | 30 |
| 2,4-Dinitrotoluene | 53 | | 65 | | 48-143 | 20 | | 30 |
| 2,6-Dinitrotoluene | 53 | | 63 | | 40-140 | 17 | | 30 |
| 4-Chlorophenyl phenyl ether | 62 | | 70 | | 40-140 | 12 | | 30 |
| 4-Bromophenyl phenyl ether | 58 | | 65 | | 40-140 | 11 | | 30 |
| Bis(2-chloroisopropyl)ether | 70 | | 79 | | 40-140 | 12 | | 30 |
| Bis(2-chloroethoxy)methane | 62 | | 70 | | 40-140 | 12 | | 30 |
| Hexachlorocyclopentadiene | 52 | | 59 | | 40-140 | 13 | | 30 |
| Isophorone | 61 | | 71 | | 40-140 | 15 | | 30 |
| Nitrobenzene | 62 | | 73 | | 40-140 | 16 | | 30 |
| NDPA/DPA | 60 | | 69 | | 40-140 | 14 | | 30 |
| n-Nitrosodi-n-propylamine | 65 | | 76 | | 29-132 | 16 | | 30 |
| Bis(2-ethylhexyl)phthalate | 64 | | 76 | | 40-140 | 17 | | 30 |
| Butyl benzyl phthalate | 54 | | 63 | | 40-140 | 15 | | 30 |
| Di-n-butylphthalate | 60 | | 71 | | 40-140 | 17 | | 30 |
| Di-n-octylphthalate | 58 | | 68 | | 40-140 | 16 | | 30 |
| Diethyl phthalate | 62 | | 70 | | 40-140 | 12 | | 30 |
| Dimethyl phthalate | 58 | | 65 | | 40-140 | 11 | | 30 |
| Biphenyl | 58 | | 64 | | 40-140 | 10 | | 30 |
| 4-Chloroaniline | 39 | Q | 67 | | 40-140 | 53 | Q | 30 |
| 2-Nitroaniline | 51 | Q | 60 | | 52-143 | 16 | | 30 |
| 3-Nitroaniline | 45 | | 44 | | 25-145 | 2 | | 30 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | LCS %Recovery | Qual | LCSD %Recovery | Qual | %Recovery Limits | RPD | Qual | RPD Limits |
|--|------------------|------|-------------------|------|---------------------|-----|------|---------------|
| Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1361431-2 WG1361431-3 | | | | | | | | |
| 4-Nitroaniline | 43 | Q | 49 | Q | 51-143 | 13 | | 30 |
| Dibenzofuran | 64 | | 70 | | 40-140 | 9 | | 30 |
| 1,2,4,5-Tetrachlorobenzene | 56 | | 62 | | 2-134 | 10 | | 30 |
| Acetophenone | 57 | | 66 | | 39-129 | 15 | | 30 |
| 2,4,6-Trichlorophenol | 55 | | 62 | | 30-130 | 12 | | 30 |
| p-Chloro-m-cresol | 59 | | 68 | | 23-97 | 14 | | 30 |
| 2-Chlorophenol | 58 | | 68 | | 27-123 | 16 | | 30 |
| 2,4-Dichlorophenol | 58 | | 66 | | 30-130 | 13 | | 30 |
| 2,4-Dimethylphenol | 53 | | 66 | | 30-130 | 22 | | 30 |
| 2-Nitrophenol | 54 | | 65 | | 30-130 | 18 | | 30 |
| 4-Nitrophenol | 50 | | 62 | | 10-80 | 21 | | 30 |
| 2,4-Dinitrophenol | 55 | | 61 | | 20-130 | 10 | | 30 |
| 4,6-Dinitro-o-cresol | 62 | | 67 | | 20-164 | 8 | | 30 |
| Phenol | 43 | | 49 | | 12-110 | 13 | | 30 |
| 2-Methylphenol | 57 | | 68 | | 30-130 | 18 | | 30 |
| 3-Methylphenol/4-Methylphenol | 57 | | 67 | | 30-130 | 16 | | 30 |
| 2,4,5-Trichlorophenol | 56 | | 63 | | 30-130 | 12 | | 30 |
| Carbazole | 62 | | 73 | | 55-144 | 16 | | 30 |
| Atrazine | 67 | | 78 | | 40-140 | 15 | | 30 |
| Benzaldehyde | 57 | | 67 | | 40-140 | 16 | | 30 |
| Caprolactam | 27 | | 31 | | 10-130 | 14 | | 30 |
| 2,3,4,6-Tetrachlorophenol | 52 | | 63 | | 40-140 | 19 | | 30 |

Lab Control Sample Analysis

Batch Quality Control

Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

| Parameter | <i>LCS</i> %Recovery | <i>Qual</i> | <i>LCSD</i> %Recovery | <i>Qual</i> | <i>%Recovery</i> Limits | <i>RPD</i> | <i>Qual</i> | <i>RPD</i> Limits |
|-----------|-------------------------|-------------|--------------------------|-------------|----------------------------|------------|-------------|----------------------|
|-----------|-------------------------|-------------|--------------------------|-------------|----------------------------|------------|-------------|----------------------|

Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1361431-2 WG1361431-3

| <i>Surrogate</i> | <i>LCS</i> %Recovery | <i>Qual</i> | <i>LCSD</i> %Recovery | <i>Qual</i> | <i>Acceptance</i> Criteria |
|----------------------|-------------------------|-------------|--------------------------|-------------|-------------------------------|
| 2-Fluorophenol | 52 | | 61 | | 21-120 |
| Phenol-d6 | 44 | | 50 | | 10-120 |
| Nitrobenzene-d5 | 59 | | 68 | | 23-120 |
| 2-Fluorobiphenyl | 55 | | 63 | | 15-120 |
| 2,4,6-Tribromophenol | 51 | | 57 | | 10-120 |
| 4-Terphenyl-d14 | 52 | | 59 | | 41-149 |

Project Name: Not Specified

Project Number: 31647

Sample Receipt and Container Information

Were project specific reporting limits specified?

YES

Cooler Information

| Cooler | Custody Seal |
|--------|--------------|
| A | Absent |

Container Information

| Container ID | Container Type | Cooler | Initial pH | Final pH | Temp deg C | Pres | Seal | Frozen Date/Time | Analysis(*) |
|--------------|-------------------------|--------|------------|----------|------------|------|--------|------------------|---|
| L2015218-01A | Vial HCl preserved | A | NA | | 5.1 | Y | Absent | | NYTCL-8260-R2(14) |
| L2015218-01B | Vial HCl preserved | A | NA | | 5.1 | Y | Absent | | NYTCL-8260-R2(14) |
| L2015218-01C | Vial HCl preserved | A | NA | | 5.1 | Y | Absent | | NYTCL-8260-R2(14) |
| L2015218-01D | Amber 250ml unpreserved | A | 7 | 7 | 5.1 | Y | Absent | | NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7) |
| L2015218-01E | Amber 250ml unpreserved | A | 7 | 7 | 5.1 | Y | Absent | | NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7) |
| L2015218-02A | Vial HCl preserved | A | NA | | 5.1 | Y | Absent | | HOLD-8260(14) |
| L2015218-02B | Vial HCl preserved | A | NA | | 5.1 | Y | Absent | | HOLD-8260(14) |

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

GLOSSARY

Acronyms

| | |
|----------|--|
| DL | - Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the limit of quantitation (LOQ). The DL includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| EDL | - Estimated Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The EDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. The use of EDLs is specific to the analysis of PAHs using Solid-Phase Microextraction (SPME). |
| EMPC | - Estimated Maximum Possible Concentration: The concentration that results from the signal present at the retention time of an analyte when the ions meet all of the identification criteria except the ion abundance ratio criteria. An EMPC is a worst-case estimate of the concentration. |
| EPA | - Environmental Protection Agency. |
| LCS | - Laboratory Control Sample: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes. |
| LCSD | - Laboratory Control Sample Duplicate: Refer to LCS. |
| LFB | - Laboratory Fortified Blank: A sample matrix, free from the analytes of interest, spiked with verified known amounts of analytes or a material containing known and verified amounts of analytes. |
| LOD | - Limit of Detection: This value represents the level to which a target analyte can reliably be detected for a specific analyte in a specific matrix by a specific method. The LOD includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| LOQ | - Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.) |
| MDL | - Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable. |
| MS | - Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values. |
| MSD | - Matrix Spike Sample Duplicate: Refer to MS. |
| NA | - Not Applicable. |
| NC | - Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit. |
| NDPA/DPA | - N-Nitrosodiphenylamine/Diphenylamine. |
| NI | - Not Ignitable. |
| NP | - Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil. |
| RL | - Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable. |
| RPD | - Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report. |
| SRM | - Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples. |
| STLP | - Semi-dynamic Tank Leaching Procedure per EPA Method 1315. |
| TEF | - Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD. |
| TEQ | - Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values. |
| TIC | - Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations. |

Footnotes

Report Format: DU Report with 'J' Qualifiers



Project Name: Not Specified

Lab Number: L2015218

Project Number: 31647

Report Date: 04/16/20

- 1 - The reference for this analyte should be considered modified since this analyte is absent from the target analyte list of the original method.

Terms

Analytical Method: Both the document from which the method originates and the analytical reference method. (Example: EPA 8260B is shown as 1,8260B.) The codes for the reference method documents are provided in the References section of the Addendum.

Difference: With respect to Total Oxidizable Precursor (TOP) Assay analysis, the difference is defined as the Post-Treatment value minus the Pre-Treatment value.

Final pH: As it pertains to Sample Receipt & Container Information section of the report, Final pH reflects pH of container determined after adjustment at the laboratory, if applicable. If no adjustment required, value reflects Initial pH.

Frozen Date/Time: With respect to Volatile Organics in soil, Frozen Date/Time reflects the date/time at which associated Reagent Water-preserved vials were initially frozen. Note: If frozen date/time is beyond 48 hours from sample collection, value will be reflected in 'bold'.

Initial pH: As it pertains to Sample Receipt & Container Information section of the report, Initial pH reflects pH of container determined upon receipt, if applicable.

PAH Total: With respect to Alkylated PAH analyses, the 'PAHs, Total' result is defined as the summation of results for all or a subset of the following compounds: Naphthalene, C1-C4 Naphthalenes, 2-Methylnaphthalene, 1-Methylnaphthalene, Biphenyl, Acenaphthylene, Acenaphthene, Fluorene, C1-C3 Fluorenes, Phenanthrene, C1-C4 Phenanthrenes/Anthracenes, Anthracene, Fluoranthene, Pyrene, C1-C4 Fluoranthenes/Pyrenes, Benz(a)anthracene, Chrysene, C1-C4 Chrysenes, Benzo(b)fluoranthene, Benzo(j)+(k)fluoranthene, Benzo(e)pyrene, Benzo(a)pyrene, Perylene, Indeno(1,2,3-cd)pyrene, Dibenz(ah)+(ac)anthracene, Benzo(g,h,i)perylene. If a 'Total' result is requested, the results of its individual components will also be reported.

PFAS Total: With respect to PFAS analyses, the 'PFAS, Total (5)' result is defined as the summation of results for: PFHpA, PFHxS, PFOA, PFNA and PFOS. If a 'Total' result is requested, the results of its individual components will also be reported.

The target compound Chlordane (CAS No. 57-74-9) is reported for GC ECD analyses. Per EPA, this compound "refers to a mixture of chlordane isomers, other chlorinated hydrocarbons and numerous other components." (Reference: USEPA Toxicological Review of Chlordane, In Support of Summary Information on the Integrated Risk Information System (IRIS), December 1997.)

Total: With respect to Organic analyses, a 'Total' result is defined as the summation of results for individual isomers or Aroclors. If a 'Total' result is requested, the results of its individual components will also be reported. This is applicable to 'Total' results for methods 8260, 8081 and 8082.

Data Qualifiers

- A** - Spectra identified as "Aldol Condensates" are byproducts of the extraction/concentration procedures when acetone is introduced in the process.
- B** - The analyte was detected above the reporting limit in the associated method blank. Flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For MCP-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank. For DOD-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte at less than ten times (10x) the concentration found in the blank AND the analyte was detected above one-half the reporting limit (or above the reporting limit for common lab contaminants) in the associated method blank. For NJ-Air-related projects, flag only applies to associated field samples that have detectable concentrations of the analyte above the reporting limit. For NJ-related projects (excluding Air), flag only applies to associated field samples that have detectable concentrations of the analyte, which was detected above the reporting limit in the associated method blank or above five times the reporting limit for common lab contaminants (Phthalates, Acetone, Methylene Chloride, 2-Butanone).
- C** - Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- D** - Concentration of analyte was quantified from diluted analysis. Flag only applies to field samples that have detectable concentrations of the analyte.
- E** - Concentration of analyte exceeds the range of the calibration curve and/or linear range of the instrument.
- G** - The concentration may be biased high due to matrix interferences (i.e. co-elution) with non-target compound(s). The result should be considered estimated.
- H** - The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
- I** - The lower value for the two columns has been reported due to obvious interference.
- J** - Estimated value. The Target analyte concentration is below the quantitation limit (RL), but above the Method Detection Limit (MDL) or Estimated Detection Limit (EDL) for SPME-related analyses. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.
- ND** - Not detected at the method detection limit (MDL) for the sample, or estimated detection limit (EDL) for SPME-related analyses.
- NJ** - Presumptive evidence of compound. This represents an estimated concentration for Tentatively Identified Compounds (TICs), where the identification is based on a mass spectral library search.
- P** - The RPD between the results for the two columns exceeds the method-specified criteria.
- Q** - The quality control sample exceeds the associated acceptance criteria. For DOD-related projects, LCS and/or Continuing Calibration

Report Format: DU Report with 'J' Qualifiers



Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

Data Qualifiers

Standard exceedences are also qualified on all associated sample results. Note: This flag is not applicable for matrix spike recoveries when the sample concentration is greater than 4x the spike added or for batch duplicate RPD when the sample concentrations are less than 5x the RL. (Metals only.)

- R** - Analytical results are from sample re-analysis.
- RE** - Analytical results are from sample re-extraction.
- S** - Analytical results are from modified screening analysis.

Project Name: Not Specified
Project Number: 31647

Lab Number: L2015218
Report Date: 04/16/20

REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - IV, 2007.

LIMITATION OF LIABILITIES

Alpha Analytical performs services with reasonable care and diligence normal to the analytical testing laboratory industry. In the event of an error, the sole and exclusive responsibility of Alpha Analytical shall be to re-perform the work at it's own expense. In no event shall Alpha Analytical be held liable for any incidental, consequential or special damages, including but not limited to, damages in any way connected with the use of, interpretation of, information or analysis provided by Alpha Analytical.

We strongly urge our clients to comply with EPA protocol regarding sample volume, preservation, cooling, containers, sampling procedures, holding time and splitting of samples in the field.



Certification Information

The following analytes are not included in our Primary NELAP Scope of Accreditation:

Westborough Facility

EPA 624/624.1: m/p-xylene, o-xylene

EPA 8260C: NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene, Azobenzene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.

EPA 8270D: NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.

SM4500: NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO₂, NO₃.

Mansfield Facility

SM 2540D: TSS

EPA 8082A: NPW: PCB: 1, 5, 31, 87, 101, 110, 141, 151, 153, 180, 183, 187.

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene, 3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

EPA TO-12 Non-methane organics

EPA 3C Fixed gases

Biological Tissue Matrix: EPA 3050B

The following analytes are included in our Massachusetts DEP Scope of Accreditation

Westborough Facility:

Drinking Water

EPA 300.0: Chloride, Nitrate-N, Fluoride, Sulfate; **EPA 353.2:** Nitrate-N, Nitrite-N; **SM4500NO3-F:** Nitrate-N, Nitrite-N; **SM4500F-C, SM4500CN-CE,**

EPA 180.1, SM2130B, SM4500CI-D, SM2320B, SM2540C, SM4500H-B, SM4500NO2-B

EPA 332: Perchlorate; **EPA 524.2:** THMs and VOCs; **EPA 504.1:** EDB, DBCP.

Microbiology: SM9215B; SM9223-P/A, SM9223B-Colilert-QT, SM9222D.

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, **EPA 350.1:** Ammonia-N, **LACHAT 10-107-06-1-B:** Ammonia-N, **EPA 351.1, SM4500NO3-F, EPA 353.2:** Nitrate-N, **SM4500P-E, SM4500P-B, E, SM4500SO4-E, SM5220D, EPA 410.4, SM5210B, SM5310C, SM4500CL-D, EPA 1664, EPA 420.1, SM4500-CN-CE, SM2540D, EPA 300:** Chloride, Sulfate, Nitrate.

EPA 624.1: Volatile Halocarbons & Aromatics,

EPA 608.3: Chlordane, Toxaphene, Aldrin, alpha-BHC, beta-BHC, gamma-BHC, delta-BHC, Dieldrin, DDD, DDE, DDT, Endosulfan I, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

EPA 625.1: SVOC (Acid/Base/Neutral Extractables), **EPA 600/4-81-045:** PCB-Oil.

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. **EPA 200.8:** Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. **EPA 245.1** Hg.

EPA 522.

Non-Potable Water

EPA 200.7: Al, Sb, As, Be, Cd, Ca, Cr, Co, Cu, Fe, Pb, Mg, Mn, Mo, Ni, K, Se, Ag, Na, Sr, TL, Ti, V, Zn.

EPA 200.8: Al, Sb, As, Be, Cd, Cr, Cu, Fe, Pb, Mn, Ni, K, Se, Ag, Na, TL, Zn.

EPA 245.1 Hg.

SM2340B

For a complete listing of analytes and methods, please contact your Alpha Project Manager.

