

**Groundwater Monitoring Report
Semi-Annual Sampling Event
First Half of 2021**

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

CHA Project Number: 031647.000

Prepared for:

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c/o Pioneer Companies
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Syracuse, New York 13202*

Prepared by:

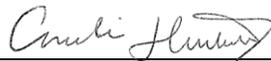


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July 2021

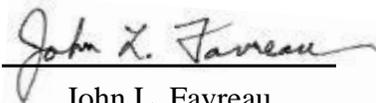
This report has been prepared and reviewed by the following qualified personnel employed by CHA.

Report Prepared By:



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Senior Scientist V

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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 2021 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 26, 2021. 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 semi volatile 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 26, 2021, 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 approximate 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 analyzed for VOCs only. Groundwater samples from the Site were historically also analyzed for metals; however, based on evaluation of historical results for detected metals, 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 2021 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 at a concentration exceeding its TOGS 1.1.1 standard; however, the detected concentration was reduced as compared to the October 2020 sampling results. No SVOCs were detected in the sample at concentrations above laboratory reporting limits.

Summary

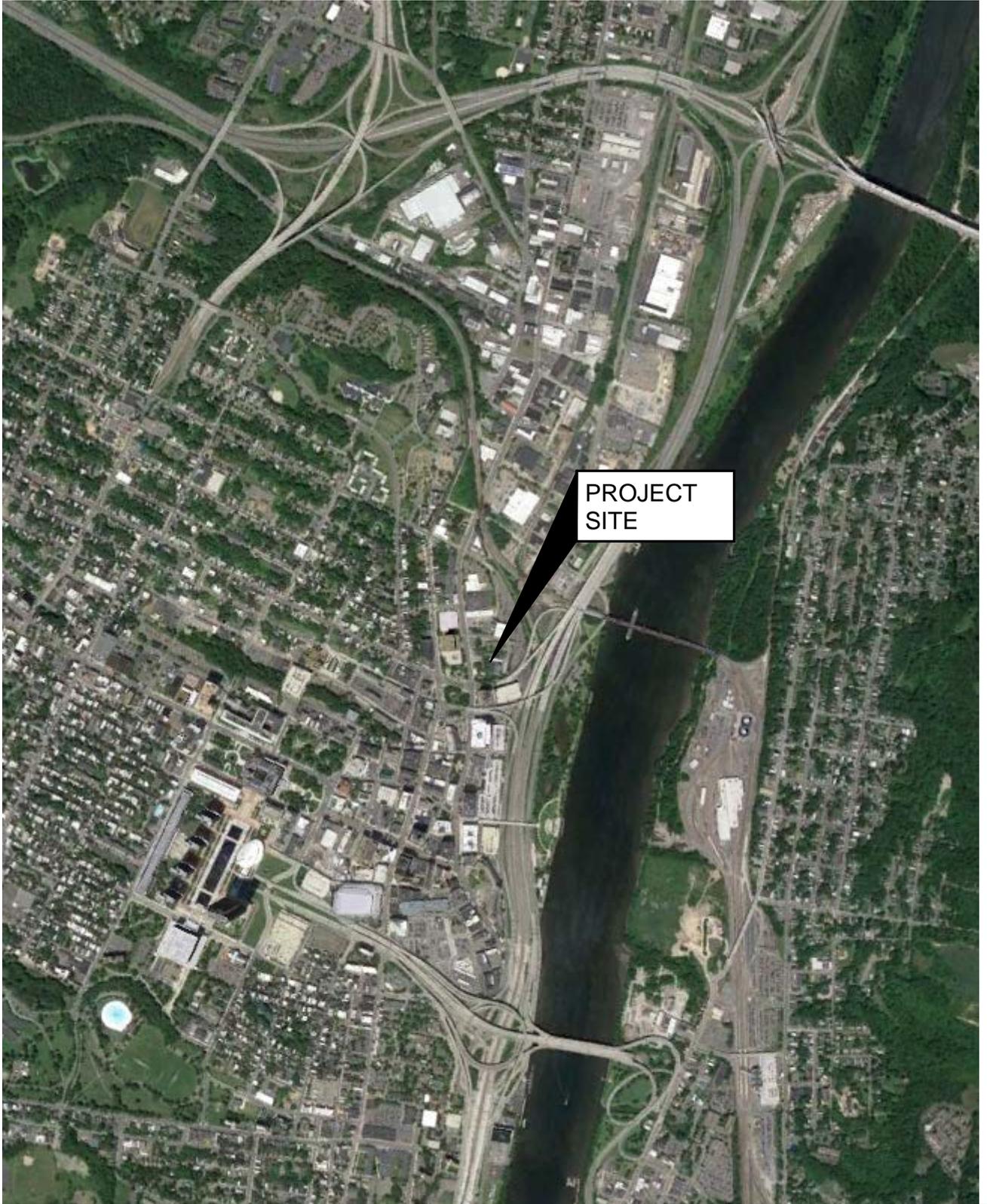
The VOC MTBE was detected at the location of well MW-1 (in the area downgradient of the former on-site gasoline station) at a concentration exceeding its TOGS 1.1.1 standard; however, the detected concentration (17 µg/l) represents a decrease from the previous sampling event in October 2020 and continuation of the decreasing trend exhibited over the past two and a half years. The detected concentration also represents a 99.5 percent decrease in the concentration of MTBE at MW-1 since September 2018. The continued decreasing trend in MTBE concentration 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 2021. Given the continued decreasing trend in MTBE concentration at well MW-1 to a concentration nearing the TOGS 1.1.1 standard, if the analytical results from the October 2021 monitoring event indicate further reduction in the concentration of MTBE, 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

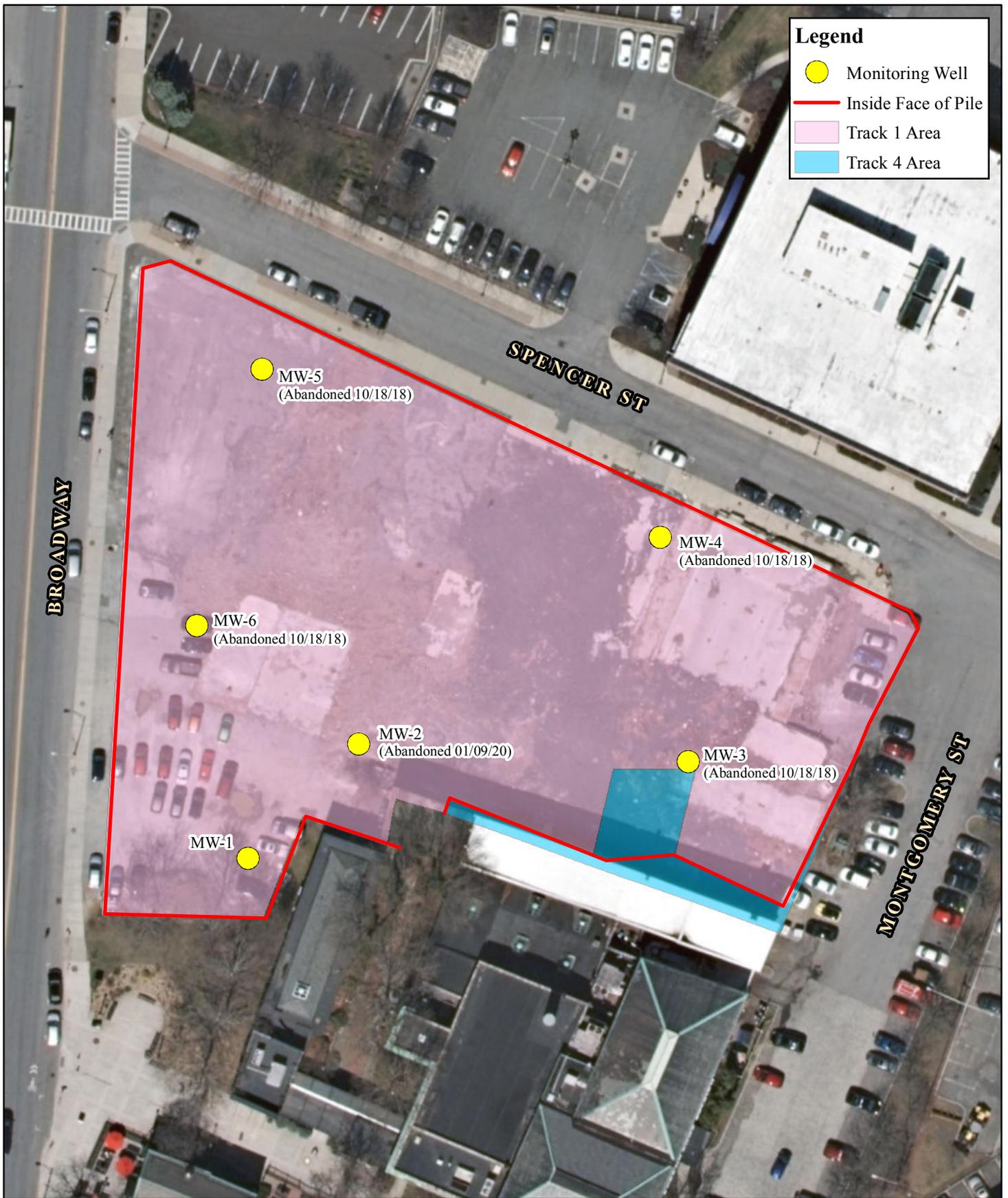


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NOT TO SCALE

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

BROADWAY

SPENCER ST

MONTGOMERY ST

MW-5
(Abandoned 10/18/18)

MW-4
(Abandoned 10/18/18)

MW-6
(Abandoned 10/18/18)

MW-2
(Abandoned 01/09/20)

MW-3
(Abandoned 10/18/18)

MW-1



Figure 2 - Site Plan
 Capital Center Site
 Quackenbush Square, 705 Broadway
 City of Albany, Albany County, New York

Scale 1" = 60'

CHA Project No.
31647

*Service Layer Credits: Aerial imagery courtesy of Esri, DigitalGlobe, GeoEye, Earthstar
Image Date: 6/7/2015*

TABLES

**Table 1
Groundwater Elevations**

Well ID		MW-1	MW-2	MW-3	MW-4	MW-5	MW-6
TOR Elevations (ft)	MW-1 DTW	103.27	101.46	96.23	94.22	101.27	101.88
3/27/2018	4.34	98.93	96.78	90.53	89.42	100.02	96.21
6/6/2018	4.85	98.42	96.17	89.98	89.61	99.52	99.21
9/27/2018	4.81	98.46	96.70	85.83	86.27	99.99	99.46
12/19/2018	6.00	97.27	96.20	decommissioned	decommissioned	decommissioned	decommissioned
3/20/2019	4.16	99.11	97.15	decommissioned	decommissioned	decommissioned	decommissioned
8/23/2019	4.54	98.73	97.84	decommissioned	decommissioned	decommissioned	decommissioned
4/9/2020	4.54	98.92	decommissioned	decommissioned	decommissioned	decommissioned	decommissioned
10/15/2020	4.54	97.98	decommissioned	decommissioned	decommissioned	decommissioned	decommissioned
4/26/2021	5.20	98.07	decommissioned	decommissioned	decommissioned	decommissioned	decommissioned

Notes:

MW-3 through MW-6 decommissioned and removed in October 2018

MW-2 decommissioned and grouted in place in January 2020

TOR - Top of Riser

TOR elevation corrected to 113.27 for 4/9/2020 and 10/15/20 events, due to raising of 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	10/15/2020	4/26/2021
			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	NA	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	NA	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	NA	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	NA	NA
Cadmium, Dissolved	5	ug/l		0.28	0.22	0.11 J	0.1 J					NA	NA	NA
Calcium, Dissolved	NS	ug/l	57600	25100	146000	166000	165000	12000	11700	7730	6720	NA	NA	NA
Chromium, Dissolved	50	ug/l					0.74 J	0.43 J	0.33 J	1.81	0.19 J	NA	NA	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	NA	NA
Copper, Dissolved	200	ug/l	1.13	0.99 J	1.25	0.93 J	2.21	1.21	1.21	3.78 J	1.39 J	NA	NA	NA
Iron, Dissolved	300	ug/l			83.2	67.9 J+	998 J+	385 J	193 J	1860 J+	293 J+	NA	NA	NA
Lead, Dissolved	25	ug/l					0.73 J			1.22		NA	NA	NA
Magnesium, Dissolved	35000	ug/l	42300	16200	103000	141000	141000 J	7290	7140	4290	3550	NA	NA	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	NA	NA
Mercury, Dissolved	0.7	ug/l			0.2 UJ							NA	NA	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	NA	NA
Potassium, Dissolved	NS	ug/l	2080	871	2490	2300	2580	822	759	1120 J	633 J	NA	NA	NA
Silver, Dissolved	50	ug/l			0.22 J							NA	NA	NA
Sodium, Dissolved	20000	ug/l	97500	73400	193000	217000	214000	68200	68100	68700	65100	NA	NA	NA
Thallium, Dissolved	0.5	ug/l									0.19 J	NA	NA	NA
Vanadium, Dissolved	NS	ug/l	2.15 J				1.85 J	2.13 J	1.82 J	6.77	4.42 J	NA	NA	NA
Zinc, Dissolved	2000	ug/l			4.41 J					8.38 J		NA	NA	NA
Total Metals														
Aluminum, Total	NS	ug/l	3790	7650	2080	16100 J	20800 J	11700	9800	12000	11900	NA	NA	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	NA	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	NA	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	NA	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	NA	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	NA	NA
Calcium, Total	NS	ug/l	75500	42400	167000	820000	844000	57200	54000	69500	78100	NA	NA	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	NA	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	NA	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	NA	NA
Iron, Total	300	ug/l	7290	14400	4980 J+	88400	107000	27700 J+	21500 J+	31600 J-	31700 J-	NA	NA	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	NA	NA
Magnesium, Total	35000	ug/l	46100	23100	113000 J+	284000	274000	21600	19500	23800	25500	NA	NA	NA
Manganese, Total	300	ug/l	417	468	1270	10080	10390	869	784.5	1053	1131	NA	NA	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	NA	NA
Potassium, Total	NS	ug/l	3230	2410	2930	6540 J	8910 J	2590	2330	2900	2820	NA	NA	NA
Selenium, Total	10	ug/l				8.83	9.54	3.37 J	3.16 J	5.71	5.33	NA	NA	NA
Silver, Total	50	ug/l		0.19 J			0.21 J					NA	NA	NA
Sodium, Total	20000	ug/l	99500	73100	200000	204000	195000	73400	74600	72900	75900	NA	NA	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	NA	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	NA	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	NA	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	22	17
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

Table 3
Summary of Purging Parameters

Well I.D.	Date	Static Water Level (ft. below TOR)	Time	Depth to Water (ft)	ORP/Eh (mV)	pH	Conductivity (mS/cm)	Turbidity (NTU)	D.O. (mg/L)	Temperature (°C)	Water Description
MW-1	4/26/2021	5.20	16:30	5.3	-180.2	7.8	0.102	40.2	4.3	21.2	Water was clear with no odor, sheen, or effervescence.
			16:35	5.32	-182.3	7.8	0.109	26.1	3.5	20.4	
			16:40	5.31	-182	7.8	0.106	10.2	3.8	20.6	
			16:45	5.3	-180.7	7.8	0.111	12.6	3.6	20.9	

APPENDIX A
MONITORING WELL SAMPLING LOG

APPENDIX B
LABORATORY ANALYTICAL DATA



ANALYTICAL REPORT

Lab Number:	L2121264
Client:	CHA Companies 3 Winners Circle Albany, NY 12205
ATTN:	C. Hurlburt
Phone:	(518) 453-4500
Project Name:	CAPITAL CENTER
Project Number:	031647.000
Report Date:	05/06/21

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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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2121264-01	MW1-04262021	WATER	ALBANY, NY	04/26/21 17:10	04/26/21
L2121264-02	TRIP BLANK	WATER	ALBANY, NY	04/26/21 00:00	04/26/21

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

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.

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:



Sebastian Corbin

Title: Technical Director/Representative

Date: 05/06/21

ORGANICS

VOLATILES

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 05/05/21 12:22
 Analyst: PD

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	ND		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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 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	17		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	ND		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	103		70-130
Toluene-d8	109		70-130
4-Bromofluorobenzene	107		70-130
Dibromofluoromethane	93		70-130

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-02
 Client ID: TRIP BLANK
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 00:00
 Date Received: 04/26/21
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8260C
 Analytical Date: 05/05/21 12:45
 Analyst: PD

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	ND		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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-02
 Client ID: TRIP BLANK
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 00:00
 Date Received: 04/26/21
 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	ND		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	ND		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	104		70-130
Toluene-d8	110		70-130
4-Bromofluorobenzene	106		70-130
Dibromofluoromethane	94		70-130

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 05/05/21 10:49
Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02 Batch: WG1494997-5					
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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 05/05/21 10:49
Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02 Batch: WG1494997-5					
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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C
Analytical Date: 05/05/21 10:49
Analyst: PD

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-02 Batch: WG1494997-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	104		70-130
Toluene-d8	106		70-130
4-Bromofluorobenzene	108		70-130
Dibromofluoromethane	94		70-130

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02 Batch: WG1494997-3 WG1494997-4								
Methylene chloride	100		100		70-130	0		20
1,1-Dichloroethane	100		100		70-130	0		20
Chloroform	99		98		70-130	1		20
Carbon tetrachloride	84		83		63-132	1		20
1,2-Dichloropropane	110		110		70-130	0		20
Dibromochloromethane	95		94		63-130	1		20
1,1,2-Trichloroethane	100		100		70-130	0		20
Tetrachloroethene	95		95		70-130	0		20
Chlorobenzene	99		99		75-130	0		20
Trichlorofluoromethane	85		86		62-150	1		20
1,2-Dichloroethane	100		100		70-130	0		20
1,1,1-Trichloroethane	92		92		67-130	0		20
Bromodichloromethane	93		95		67-130	2		20
trans-1,3-Dichloropropene	98		97		70-130	1		20
cis-1,3-Dichloropropene	96		94		70-130	2		20
Bromoform	96		96		54-136	0		20
1,1,2,2-Tetrachloroethane	110		110		67-130	0		20
Benzene	98		97		70-130	1		20
Toluene	100		100		70-130	0		20
Ethylbenzene	98		97		70-130	1		20
Chloromethane	81		80		64-130	1		20
Bromomethane	51		57		39-139	11		20
Vinyl chloride	90		90		55-140	0		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02 Batch: WG1494997-3 WG1494997-4								
Chloroethane	100		98		55-138	2		20
1,1-Dichloroethene	91		91		61-145	0		20
trans-1,2-Dichloroethene	98		97		70-130	1		20
Trichloroethene	96		96		70-130	0		20
1,2-Dichlorobenzene	100		100		70-130	0		20
1,3-Dichlorobenzene	100		100		70-130	0		20
1,4-Dichlorobenzene	100		100		70-130	0		20
Methyl tert butyl ether	95		96		63-130	1		20
p/m-Xylene	100		95		70-130	5		20
o-Xylene	100		100		70-130	0		20
cis-1,2-Dichloroethene	100		100		70-130	0		20
Styrene	95		100		70-130	5		20
Dichlorodifluoromethane	63		64		36-147	2		20
Acetone	140		120		58-148	15		20
Carbon disulfide	97		94		51-130	3		20
2-Butanone	110		110		63-138	0		20
4-Methyl-2-pentanone	110		110		59-130	0		20
2-Hexanone	110		110		57-130	0		20
Bromochloromethane	100		100		70-130	0		20
1,2-Dibromoethane	100		99		70-130	1		20
1,2-Dibromo-3-chloropropane	88		89		41-144	1		20
Isopropylbenzene	110		110		70-130	0		20
1,2,3-Trichlorobenzene	96		100		70-130	4		20

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02 Batch: WG1494997-3 WG1494997-4								
1,2,4-Trichlorobenzene	100		100		70-130	0		20
Methyl Acetate	96		100		70-130	4		20
Cyclohexane	97		95		70-130	2		20
1,4-Dioxane	80		88		56-162	10		20
Freon-113	84		84		70-130	0		20
Methyl cyclohexane	86		85		70-130	1		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	101		104		70-130
Toluene-d8	106		106		70-130
4-Bromofluorobenzene	108		108		70-130
Dibromofluoromethane	99		97		70-130

SEMIVOLATILES

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8270D
 Analytical Date: 05/04/21 01:11
 Analyst: EK

Extraction Method: EPA 3510C
 Extraction Date: 05/02/21 14:34

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	1.6	J	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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 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	49		21-120
Phenol-d6	39		10-120
Nitrobenzene-d5	57		23-120
2-Fluorobiphenyl	67		15-120
2,4,6-Tribromophenol	85		10-120
4-Terphenyl-d14	69		41-149

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 Field Prep: Not Specified

Sample Depth:

Matrix: Water
 Analytical Method: 1,8270D-SIM
 Analytical Date: 05/06/21 12:50
 Analyst: DV

Extraction Method: EPA 3510C
 Extraction Date: 05/02/21 14:36

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	0.02	J	ug/l	0.10	0.01	1
Phenanthrene	0.03	J	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	0.06	J	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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

SAMPLE RESULTS

Lab ID: L2121264-01
 Client ID: MW1-04262021
 Sample Location: ALBANY, NY

Date Collected: 04/26/21 17:10
 Date Received: 04/26/21
 Field Prep: Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
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Semivolatile Organics by GC/MS-SIM - Westborough Lab

Surrogate	% Recovery	Qualifier	Acceptance Criteria
2-Fluorophenol	48		21-120
Phenol-d6	43		10-120
Nitrobenzene-d5	66		23-120
2-Fluorobiphenyl	71		15-120
2,4,6-Tribromophenol	59		10-120
4-Terphenyl-d14	85		41-149

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D
Analytical Date: 05/04/21 00:22
Analyst: EK

Extraction Method: EPA 3510C
Extraction Date: 05/02/21 14:34

Parameter	Result	Qualifier	Units	RL	MDL
Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1493452-1					
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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

**Method Blank Analysis
Batch Quality Control**

Analytical Method: 1,8270D
Analytical Date: 05/04/21 00:22
Analyst: EK

Extraction Method: EPA 3510C
Extraction Date: 05/02/21 14:34

Parameter	Result	Qualifier	Units	RL	MDL
Semivolatile Organics by GC/MS - Westborough Lab for sample(s): 01 Batch: WG1493452-1					
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	72		21-120
Phenol-d6	59		10-120
Nitrobenzene-d5	112		23-120
2-Fluorobiphenyl	97		15-120
2,4,6-Tribromophenol	77		10-120
4-Terphenyl-d14	101		41-149

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D-SIM
Analytical Date: 05/04/21 13:33
Analyst: JRW

Extraction Method: EPA 3510C
Extraction Date: 05/02/21 14:36

Parameter	Result	Qualifier	Units	RL	MDL
Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01 Batch: WG1493454-1					
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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8270D-SIM
Analytical Date: 05/04/21 13:33
Analyst: JRW

Extraction Method: EPA 3510C
Extraction Date: 05/02/21 14:36

Parameter	Result	Qualifier	Units	RL	MDL
Semivolatile Organics by GC/MS-SIM - Westborough Lab for sample(s): 01 Batch: WG1493454-1					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
2-Fluorophenol	64		21-120
Phenol-d6	55		10-120
Nitrobenzene-d5	89		23-120
2-Fluorobiphenyl	98		15-120
2,4,6-Tribromophenol	78		10-120
4-Terphenyl-d14	125		41-149

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Lab Number: L2121264

Project Number: 031647.000

Report Date: 05/06/21

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: WG1493452-2 WG1493452-3								
Bis(2-chloroethyl)ether	65		65		40-140	0		30
3,3'-Dichlorobenzidine	64		63		40-140	2		30
2,4-Dinitrotoluene	73		76		48-143	4		30
2,6-Dinitrotoluene	79		79		40-140	0		30
4-Chlorophenyl phenyl ether	62		64		40-140	3		30
4-Bromophenyl phenyl ether	63		62		40-140	2		30
Bis(2-chloroisopropyl)ether	81		80		40-140	1		30
Bis(2-chloroethoxy)methane	77		72		40-140	7		30
Hexachlorocyclopentadiene	61		61		40-140	0		30
Isophorone	80		77		40-140	4		30
Nitrobenzene	84		81		40-140	4		30
NDPA/DPA	68		70		40-140	3		30
n-Nitrosodi-n-propylamine	88		82		29-132	7		30
Bis(2-ethylhexyl)phthalate	90		92		40-140	2		30
Butyl benzyl phthalate	88		87		40-140	1		30
Di-n-butylphthalate	78		77		40-140	1		30
Di-n-octylphthalate	95		98		40-140	3		30
Diethyl phthalate	73		74		40-140	1		30
Dimethyl phthalate	77		73		40-140	5		30
Biphenyl	70		66		40-140	6		30
4-Chloroaniline	75		78		40-140	4		30
2-Nitroaniline	82		81		52-143	1		30
3-Nitroaniline	63		68		25-145	8		30

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

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: WG1493452-2 WG1493452-3								
4-Nitroaniline	65		65		51-143	0		30
Dibenzofuran	64		65		40-140	2		30
1,2,4,5-Tetrachlorobenzene	66		64		2-134	3		30
Acetophenone	71		71		39-129	0		30
2,4,6-Trichlorophenol	75		68		30-130	10		30
p-Chloro-m-cresol	89		82		23-97	8		30
2-Chlorophenol	70		66		27-123	6		30
2,4-Dichlorophenol	74		74		30-130	0		30
2,4-Dimethylphenol	75		76		30-130	1		30
2-Nitrophenol	90		91		30-130	1		30
4-Nitrophenol	54		57		10-80	5		30
2,4-Dinitrophenol	65		61		20-130	6		30
4,6-Dinitro-o-cresol	69		73		20-164	6		30
Phenol	56		55		12-110	2		30
2-Methylphenol	68		70		30-130	3		30
3-Methylphenol/4-Methylphenol	72		72		30-130	0		30
2,4,5-Trichlorophenol	81		67		30-130	19		30
Carbazole	76		75		55-144	1		30
Atrazine	90		89		40-140	1		30
Benzaldehyde	65		61		40-140	6		30
Caprolactam	47		43		10-130	9		30
2,3,4,6-Tetrachlorophenol	65		68		40-140	5		30

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
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Semivolatile Organics by GC/MS - Westborough Lab Associated sample(s): 01 Batch: WG1493452-2 WG1493452-3

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
2-Fluorophenol	76		73		21-120
Phenol-d6	60		59		10-120
Nitrobenzene-d5	106		106		23-120
2-Fluorobiphenyl	90		88		15-120
2,4,6-Tribromophenol	81		90		10-120
4-Terphenyl-d14	90		99		41-149

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

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: WG1493454-2 WG1493454-3								
Acenaphthene	80		86		40-140	7		40
2-Chloronaphthalene	78		84		40-140	7		40
Fluoranthene	91		103		40-140	12		40
Hexachlorobutadiene	66		72		40-140	9		40
Naphthalene	74		80		40-140	8		40
Benzo(a)anthracene	83		89		40-140	7		40
Benzo(a)pyrene	90		98		40-140	9		40
Benzo(b)fluoranthene	90		103		40-140	13		40
Benzo(k)fluoranthene	92		97		40-140	5		40
Chrysene	83		93		40-140	11		40
Acenaphthylene	82		89		40-140	8		40
Anthracene	85		94		40-140	10		40
Benzo(ghi)perylene	85		91		40-140	7		40
Fluorene	85		91		40-140	7		40
Phenanthrene	82		90		40-140	9		40
Dibenzo(a,h)anthracene	92		99		40-140	7		40
Indeno(1,2,3-cd)pyrene	89		93		40-140	4		40
Pyrene	90		102		40-140	13		40
2-Methylnaphthalene	81		87		40-140	7		40
Pentachlorophenol	77		84		40-140	9		40
Hexachlorobenzene	73		79		40-140	8		40
Hexachloroethane	64		70		40-140	9		40

Lab Control Sample Analysis

Batch Quality Control

Project Name: CAPITAL CENTER

Project Number: 031647.000

Lab Number: L2121264

Report Date: 05/06/21

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
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Semivolatile Organics by GC/MS-SIM - Westborough Lab Associated sample(s): 01 Batch: WG1493454-2 WG1493454-3

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
2-Fluorophenol	70		76		21-120
Phenol-d6	60		63		10-120
Nitrobenzene-d5	89		97		23-120
2-Fluorobiphenyl	95		103		15-120
2,4,6-Tribromophenol	88		103		10-120
4-Terphenyl-d14	123		140		41-149

Project Name: CAPITAL CENTER**Lab Number:** L2121264**Project Number:** 031647.000**Report Date:** 05/06/21**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(*)
L2121264-01A	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2121264-01B	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2121264-01C	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2121264-01D	Amber 250ml unpreserved	A	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2121264-01E	Amber 250ml unpreserved	A	7	7	3.4	Y	Absent		NYTCL-8270-SIM-LVI(7),NYTCL-8270-LVI(7)
L2121264-02A	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)
L2121264-02B	Vial HCl preserved	A	NA		3.4	Y	Absent		NYTCL-8260-R2(14)

Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

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.
NR	- No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.
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.

Report Format: DU Report with 'J' Qualifiers



Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Footnotes

- 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. In addition, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. (Note: 'PFAS, Total (6)' is applicable to MassDEP DW compliance analysis only.). 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.
- F** - The ratio of quantifier ion response to qualifier ion response falls outside of the laboratory criteria. Results are considered to be an estimated maximum concentration.
- 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.

Report Format: DU Report with 'J' Qualifiers



Project Name: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

Data Qualifiers

- 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 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: CAPITAL CENTER
Project Number: 031647.000

Lab Number: L2121264
Report Date: 05/06/21

REFERENCES

- 1 Test Methods for Evaluating Solid Waste: Physical/Chemical Methods. EPA SW-846. Third Edition. Updates I - VI, 2018.

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, Naphthalene

EPA 625/625.1: alpha-Terpineol

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

EPA 8270D/8270E: NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol; 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.

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, **LCHAT 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, SM9222D.**

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, EPA 537.1.

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.

