

November 12, 2024

Megan Kuczka, DER Project Manager  
New York State Department of Environmental Conservation  
Division of Environmental Remediation, Region 9  
700 Delaware Avenue  
Buffalo, New York 14209

Re: **Monitoring and Sampling Summary (3rd Quarter 2024)**  
Site Management Plan, Post Installation Monitoring & Inspection  
MOD-PAC CORP. Site, 1801 Elmwood Avenue, Buffalo, New York

Dear Ms. Kuczka:

In accordance with the Site Management Plan (SMP)<sup>1</sup> for NYSDEC Site #C915314, Environmental Advantage, Inc. (EA), has prepared this summary letter report which provides the results of the inspection, monitoring, and maintenance of the Sub-Slab Depressurization (SSD) systems completed from July 1 through September 30, 2024. The attachments to this letter report include figures (Attachment A), summary tables (Attachment B), well data sheets (Attachment C), and analytical laboratory reports (Attachment D).

After discussions with the New York State Department of Environmental Conservation (NYSDEC or Department), New York State Department of Health (NYSDOH) representatives, and Matrix Environmental Technologies, Inc. (METI), the engineering firm responsible for the design and annual inspection and certification of the SSD systems, it was determined that quarterly groundwater sampling of the Site's four groundwater monitoring wells subject to the remedial program was warranted to investigate seasonal variation in contaminant concentrations and the potential seasonal correlation to maintaining a negative pressure of at least 0.002 inches water column (WC) in the sub-slab as the SSD Systems were designed. The monthly collection of vacuum readings for any vapor monitoring point (VMP) that fails to achieve the minimum negative pressure of at least 0.002 inches WC during quarterly SSD inspections was also initiated, until the affected VMP('s) meet the minimum negative pressure as designed (with the exception of VMP-6A<sup>2</sup>, VMP-8A, and VMP-5B which are considered inactive). In April 2024, VMP-8A was replaced with VMP-8AR, and VMP-5B was replaced with VMP-5BR due to VMP-8A and VMP-5B frequently failing to achieve the minimum negative pressure of at least 0.002 inches WC. The locations of the groundwater monitoring wells, and SSD systems are shown on Figure 1.

<sup>1</sup> "Site Management Plan for MOD-PAC Site, 1801 Elmwood Avenue, City of Buffalo, Erie County, New York, Site No. C915314" prepared by C&S Engineers, Inc., December 2019, revised March 2022 by Environmental Advantage, Inc.

<sup>2</sup> VMP-6A has been verified as a dead point, as described in Section 5.1 – 'Area A Testing' of METI's "System Start-up Report and Operation and Maintenance Plan"<sup>2</sup> as provided within Appendix H – Operation and Maintenance Manual of the SMP. VMP-6A always exhibits positive pressure readings.

## **Post-Installation SSD Maintenance and Monitoring**

System checks are completed on a quarterly basis, at a minimum. Routine monitoring includes the identification and repair of any leaks, operational status checks of blowers and fans, documentation of manifold settings and vacuum point at each vapor extraction point, and documentation of vacuum at each monitoring point. During the quarterly system checks, pre- and post-carbon air samples are collected from Area A. Samples are submitted for laboratory analysis of volatile organic compounds (VOCs) via Environmental Protection Agency (EPA) Method TO-15. In addition, pre- and post-carbon photoionization detector (PID) readings are collected from Area A, as well as from Areas B and C effluent, on a monthly basis. Non-routine maintenance, including carbon change outs, is completed as necessary based on analytical data of pre- and post-carbon samples.

SSD System layout for each area is shown on Figure 2A for Area A, Figure 2B for Area B, and Figure 2C for Area C, presented in Attachment A. Area-specific findings during Q3 2024 monitoring event are summarized in Table 1 with historical data presented in Table 2A for Area A, Table 2B for Area B, and Table 2C for Area C, all of which are provided in Attachment B. Air sample results for the current monitoring period are summarized in Table 3.

### **SSD Area A – Finished Product Storage Area**

During Q3 2024, manometer readings for all active VMPs in Area A achieved the minimum negative pressure of at least 0.002 inches WC in the sub-slab with the exception of VMP-6A (dead point) in September.

Post-carbon analytical data exhibited lower concentrations of all target chlorinated compounds when compared to pre-carbon concentrations, with the exception of methylene chloride, which is inconclusive due to sample dilution. **Please Note:** the reporting limit for methylene chloride was 4.12 ug/m<sup>3</sup> for the pre-carbon sample, therefore it cannot be determined if the pre-carbon value of non-detect is higher than post-carbon concentration of 3.65 ug/m<sup>3</sup> for this analyte. Overall target chlorinated VOC (cVOC)<sup>3</sup> reduction of 99.1 percent. Air sample results for Q3 2024 are summarized in Table 3, with historical air sample results summarized in Table 4. The complete analytical laboratory report is provided in Attachment C.

### **SSD Area B – Roll Storage Area (Formerly Cold Storage Area)**

During Q3 2024, manometer readings for all active VMPs achieved the minimum 0.002 inches WC in the sub-slab with the exception of VMP-5B in September. VMP-5BR, located approximately +/- 8' to the east of VMP-5B, achieved a minimum negative pressure of at least 0.002 inches WC in July and September.

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<sup>3</sup> NYSDOH Target cVOCs are included in this calculation, specifically those listed in the NYSDOH "Final Guidance for Evaluating Soil Vapor Intrusion in the State of New York", May 2017 Update. Specifically: 1,1,1-Trichloroethane, 1,1-Dichloroethene, Carbon tetrachloride, cis-1,2-Dichloroethene, Methylene chloride, Tetrachloroethene, Trichloroethene, and Vinyl chloride

### **SSD Area C – Maintenance Area**

During Q3 2024, manometer readings for all active VMPs achieved the minimum 0.002 inches WC in the sub-slab.

### **Groundwater Monitoring**

During the current monitoring period, water table measurements were collected in July. All six wells in the vicinity of SSDS Area A, Area B, and Area C (MW-3, MW-11, MW-12, MW-13, MW-14, and MW-15) were gauged. Groundwater samples were collected on July 1, 2024 from the four monitoring wells included in the remedial program: MW-3, MW-11, MW-12, and MW-13. All samples were submitted for laboratory analysis of Target Compound List (TCL) VOCs via EPA Method 8260. Historical water table measurements for the six wells in the vicinity of SSDS Area A, Area B, and Area C are summarized in Table 5. Historical groundwater elevation monitoring and sampling data results of four monitoring wells included in the remedial program are summarized in Table 6. The complete analytical laboratory report is provided in Attachment D. **Please Note:** Groundwater elevation data are available for the four monitoring wells included in the remedial program only, the well details on MW-14 and MW-15 are not included in the Site's remedial documents.

### **Corrective Measures**

The granular activated carbon was replaced from the carbon vessel in Area A on July 18, 2024. The spent carbon was placed in 55-gallon drums and is currently pending transport for regeneration. No other additional corrective actions occurred during the Q3 2024 monitoring period.

### **Conclusions and Scheduling**

During the Q3 2024 monitoring period, all active manometers met the minimum 0.002 inches WC in the sub-slab with the exception of VMP-6A (dead point) and VMP-5B (redrilled) in September. Replacement VMP's VMP-8AR and VMP-5BR were installed in the vicinity of VMP-8A and VMP-5B, respectively, due to periodic failure to obtain the minimum negative pressure of at least 0.002 inches WC. Compliant vacuum readings were obtained at both replacement VMP's during the quarterly monitoring period. The granular activated carbon was replaced from the carbon vessel in Area A on July 18, 2024. There are no additional corrective actions to report for the Q3 2024 monitoring period. The SSD systems in Area A, Area B, and Area C appear to be functioning properly.

Post-carbon analytical data collected during Q3 2024 exhibited lower concentrations of all target chlorinated compounds, with the exception of methylene chloride, which is inconclusive. A few non-chlorinated compounds exhibited higher concentrations post-carbon compared to pre-carbon; however, the removal efficiency of carbon is much reduced at lower influent concentrations, which is likely playing a role in many of the low-level increases occasionally observed at this Site. Overall target chlorinated VOC (cVOC) was 99.1 percent. Continued system inspections, monitoring, and sampling will be completed for the fourth quarter of 2024.

If you have any questions regarding the information presented above, please contact me directly for further information.

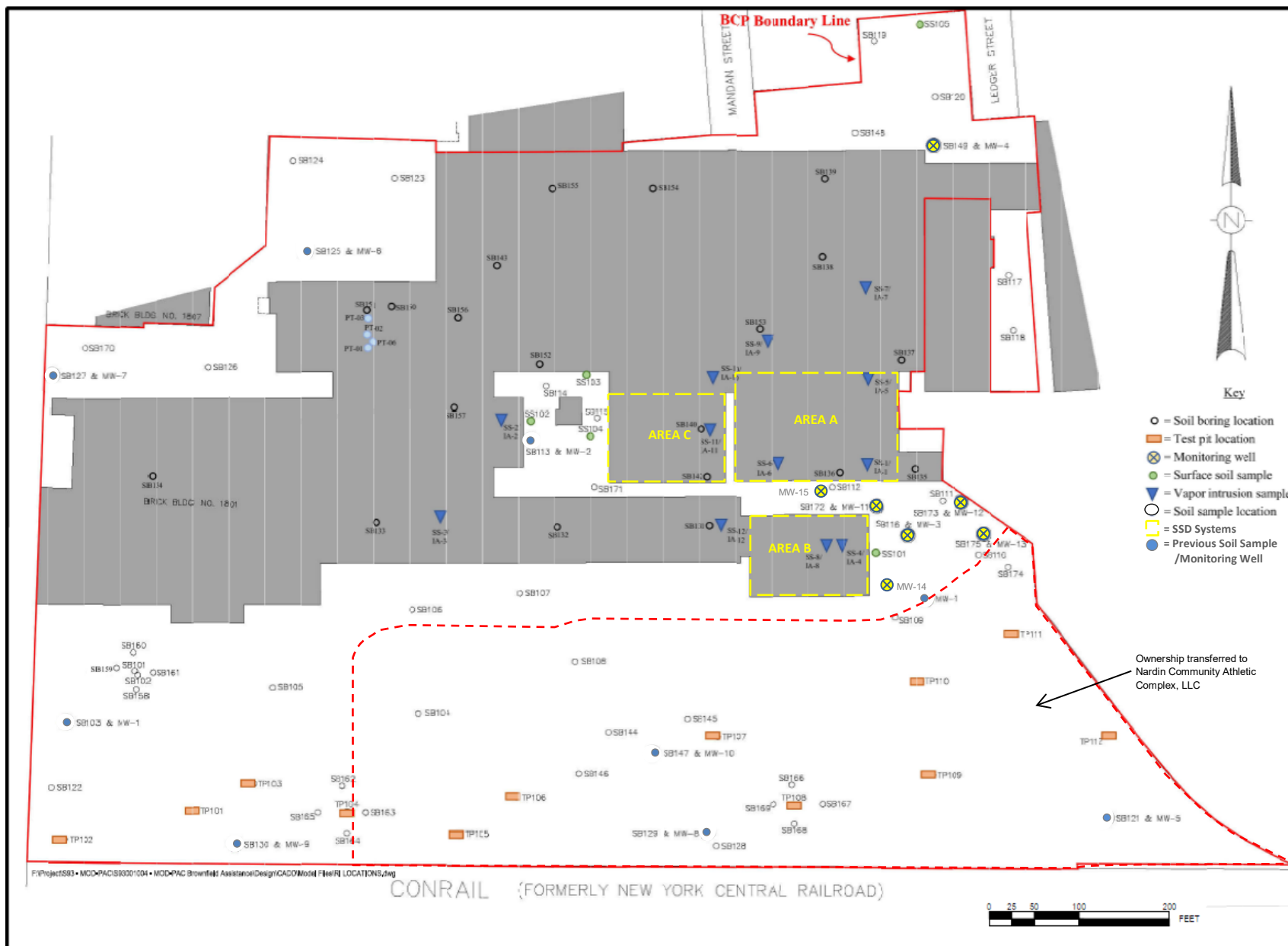
Very truly yours,  
ENVIRONMENTAL ADVANTAGE, INC.

A handwritten signature in blue ink, appearing to read "C. Mark Hanna".

C. Mark Hanna, CHMM  
President

## **ATTACHMENT A**

### Figures



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## ENVIRONMENTAL ADVANTAGE, INC.

Regulatory Compliance – Site Investigations – Facility Inspections

### BCP SITE PLAN

### MOD-PAC, CORP.

1801 ELMWOOD AVENUE  
BUFFALO, NEW YORK

DRAWN BY: MB

SCALE: NOT TO SCALE

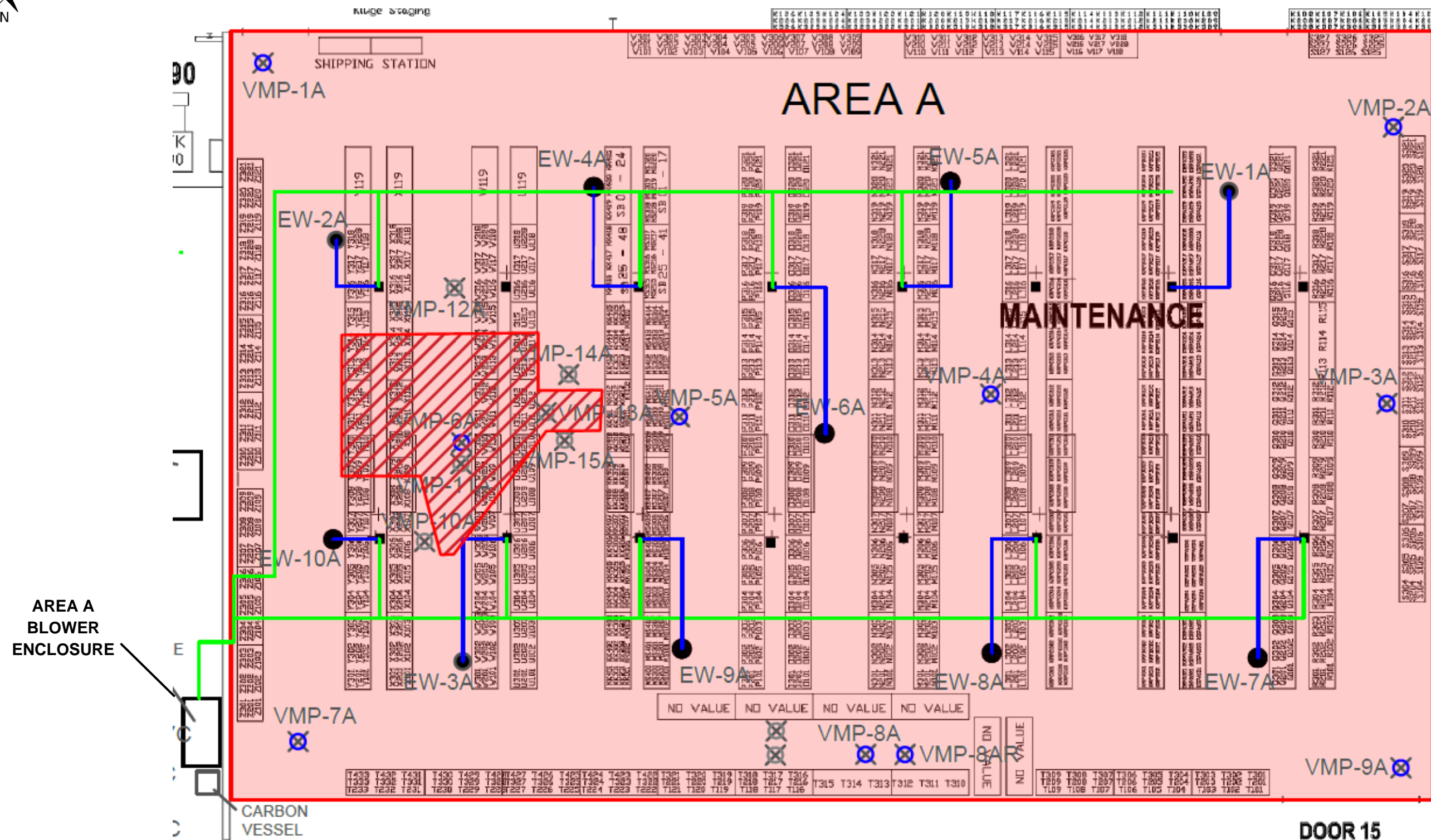
PROJECT: 01304

CHECKED BY: CMH

DATE: 06/2023

FIGURE NO: 1



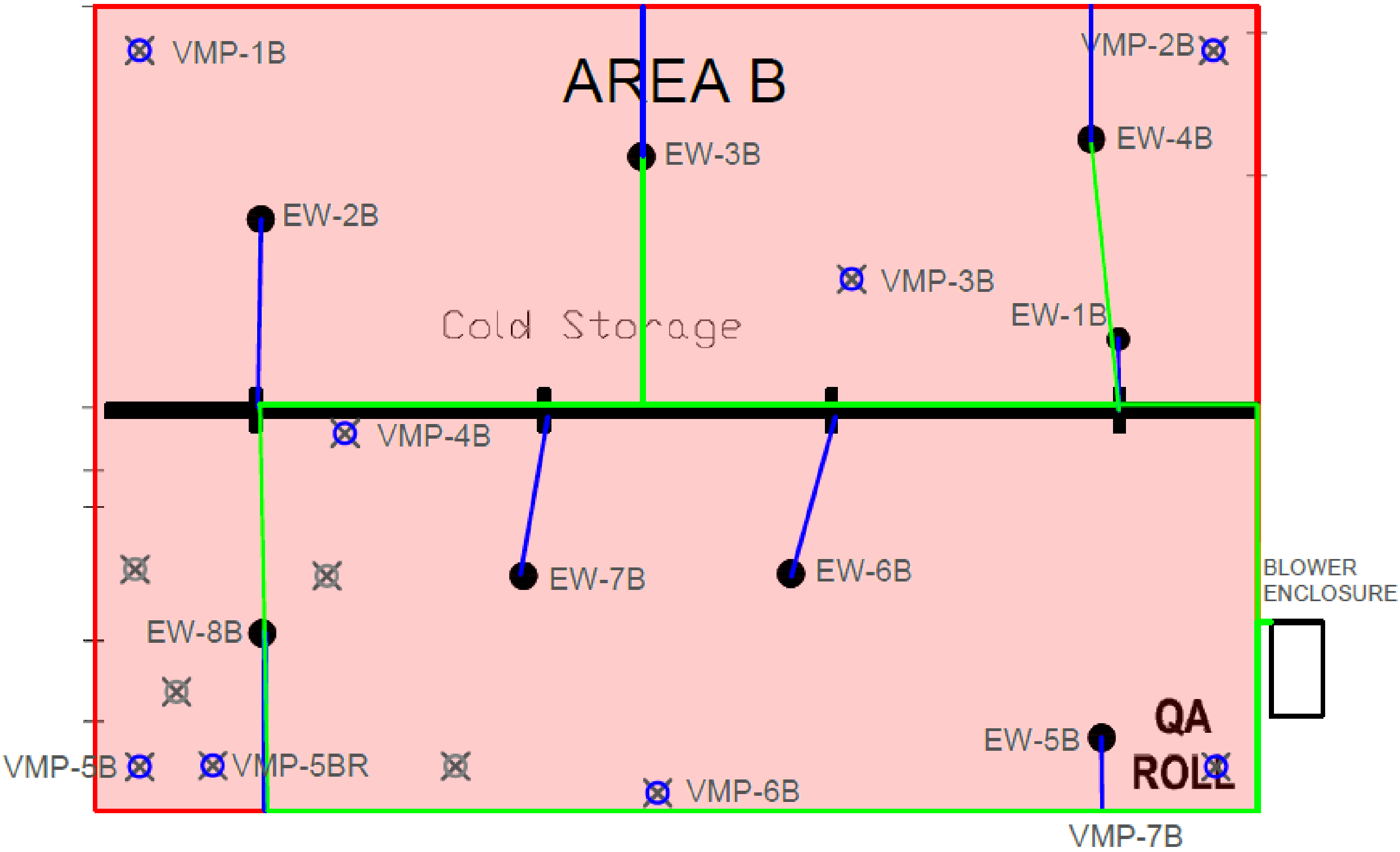


- Vapor Monitoring Point (permanent)
- Vapor Monitoring Point (temporary/decommissioned)
- Vapor Extraction Well
- Piping below grade
- Piping above grade
- Area of Zero Vacuum Influence

ENVIRONMENTAL ADVANTAGE, INC.		
Regulatory Compliance – Site Investigations – Facility Inspections		
<b>SSDS AREA A</b> 1801 ELMWOOD AVENUE BUFFALO, NEW YORK		
DRAWN BY: KL	SCALE: NOT TO SCALE	PROJECT: 01304
CHECKED BY: MS	DATE: 08/2024	FIGURE NO: 2A



THIS FIGURE WAS ADAPTED FROM SUB-SLAB DEPRESSURIZATION SYSTEMS 2024 PERIODIC REVIEW REPORT PREPARED BY MATRIX FOR MOD-PAC CORPORATION (AUGUST 2024)



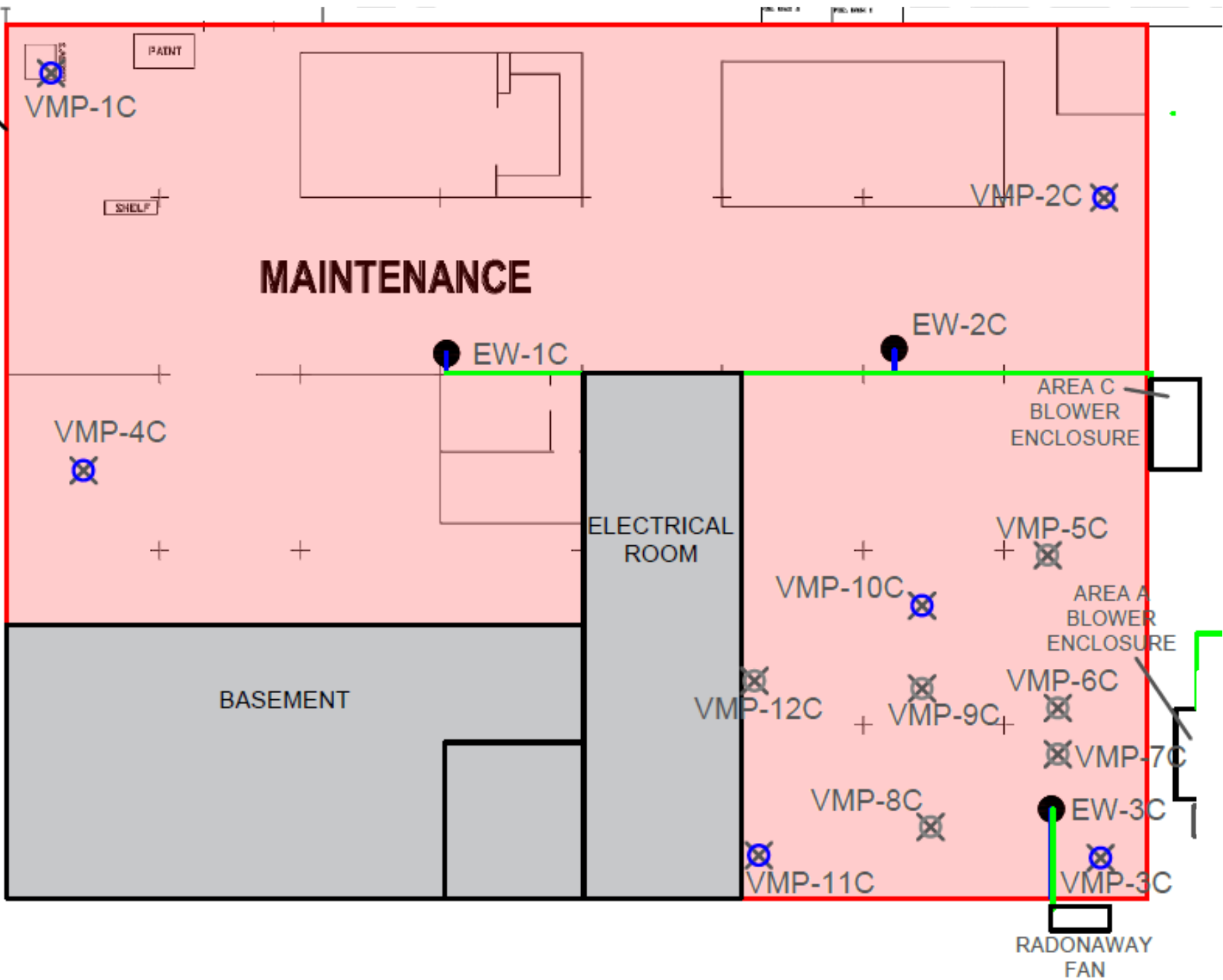
- ⊗ Vapor Monitoring Point (permanent)
- ⊗ Vapor Monitoring Point (temporary/decommissioned)
- Vapor Extraction Well
- Piping below grade
- Piping above grade
- ▨ Area of Zero Vacuum Influence

ENVIRONMENTAL ADVANTAGE, INC.		
Phase I/II Audits – Site Investigations – Facility Inspections		
SSDS AREA B		
1801 ELMWOOD AVENUE BUFFALO, NEW YORK		
DRAWN BY: KL	SCALE: NOT TO SCALE	PROJECT: 01304
CHECKED BY: MS	DATE: 08/2024	FIGURE NO: 2B





# AREA C



- Vapor Monitoring Point (permanent)
- Vapor Monitoring Point (temporary/decommissioned)
- Vapor Extraction Well
- Piping below grade
- Piping above grade
- Area of Zero Vacuum Influence

ENVIRONMENTAL ADVANTAGE, INC.		
Phase I/II Audits – Site Investigations – Facility Inspections		
SSDS AREA C		
1801 ELMWOOD AVENUE		
BUFFALO, NEW YORK		
DRAWN BY: KL	SCALE: NOT TO SCALE	PROJECT: 01304
CHECKED BY: MS	DATE: 08/2024	FIGURE NO: 2C

## **ATTACHMENT B**

### Tables

Table 1  
MOD-PAC CORP., 1801 Elmwood Ave, Buffalo, NY  
SSDS Post Installation Monitoring Results  
September Q3 2024 Summary

**Area A - Finished Product Storage Area**

Date	Extraction Wells (in WC)										Blower (in WC)	Pre-carbon PID Reading (ppm)	Post-carbon PID Reading (ppm)
	EW-1A	EW-2A	EW-3A	EW-4A	EW-5A	EW-6A	EW-7A	EW-8A	EW-9A	EW-10A			
9/6/2024	18.0	19.0	19.0	19.0	18.0	0.0	18.0	20.0	19.0	19.0	21	1.3	0.2

Date	Vapor Monitoring Points (in WC)									
	VMP-1A	VMP-2A	VMP-3A	VMP-4A	VMP-5A	VMP-6A	VMP-7A	VMP-8A	VMP-8AR	VMP-9A
9/6/2024	-0.116	-0.065	-0.093	-0.154	-0.065	+0.000	-0.059	-0.034	-0.046	-0.130

**Area B - Cold Storage Garage**

Date	Extraction Wells (in WC)								Blower (in WC)	System Effluent PID Reading (ppm)
	EW-1B	EW-2B	EW-3B	EW-4B	EW-5B	EW-6B	EW-7B	EW-8B		
9/6/2024	39.0	39.0	40.0	40.0	39.0	40.0	39.0	39.0	39.0	0.0

Date	Vapor Monitoring Points (in WC)							
	VMP-1B	VMP-2B	VMP-3B	VMP-4B	VMP-5B	VMP-5BR	VMP-6B	VMP-7B
9/6/2024	-0.018	-0.052	-0.350	-0.373	+ 0.000	-0.029	-0.015	-0.297

**Area C - Maintenance Area**

Date	Extraction Wells (in WC)			System Effluent PID Reading (ppm)		
	EW-1C	EW-2C	EW-3C	EW-1C	EW-2C	EW-3C
9/6/2024	42.0	45.0	30.0	0.7	0.8	0.3

Date	Vapor Monitoring Points (in WC)					
	VMP-1C	VMP-2C	VMP-3C	VMP-4C	VMP-10C	VMP-11C
9/6/2024	-0.051	-0.094	-0.023	-0.073	-0.124	-0.047

**Note:**

1. in WC = inches water column; ppm = parts per million;

**Table 2A**  
**MOD-PAC CORP., 1801 Elmwood Ave, Buffalo, NY**  
**SSDS Post Installation Monitoring Results**  
**Area A - Finished Product Storage Area**

Date	Extraction Wells (in WC)										Blower (in WC)	Pre-carbon PID Reading (ppm)	Post-carbon PID Reading (ppm)
	EW-1A	EW-2A	EW-3A	EW-4A	EW-5A	EW-6A	EW-7A	EW-8A	EW-9A	EW-10A			
9/26/2019	14.5	14.5	15.5	14.5	15	14.5	15	14.5	14.5	15.5	12	3.3	1.5
10/3/2019	14	14	15	14	14	1	14	15	14	15	12	52.6	12.7
10/9/2019	13	13.5	14	13.5	13.5	1	13.5	14	13.5	14.5	13	0.0	0.0
11/5/2019	11.5	12	12.5	11.5	12	1	12	12	11.5	12.5	10	4.7	0.5
12/3/2019	11	11.5	12	11	11.5	1	11.5	11.5	11.5	12	10	1.0	0.1
1/22/2020												0.2	0.0
2/11/2020	10	10.5	11	10.5	11	1	11	11	10.5	11.5	9	0.5	0.0
3/27/2020	10	10	11	10.5	11	1	10.5	10.5	10	11	8	47.8	27.1
6/28/2020	13	13	13.5	13	13	1	13	13	13	13.5	14	0.4	0.0
7/31/2020												0.0	0.0
8/28/2020												0.0	0.0
9/15/2020	13.5	14.0	14.5	14.0	14.0	1.0	14.0	14.5	14.5	15.0	14	2.7	1.1
10/19/2020												7.8	4.6
11/4/2020												0.0	0.0
12/8/2020	12.5	13.0	13.5	13.0	13.0	1.0	13.0	14.0	13.0	14.0	12	0.6	0.0
1/4/2021												0.4	0.0
2/18/2021												1.0	0.0
3/30/2021	13.0	14.0	14.0	14.0	14.0	0.0	14.0	14.0	14.0	15.0	12	0.0	0.0
4/14/2021												0.4	0.0
5/20/2021												0.4	0.0
6/11/2021	16.0	16.0	16.0	16.0	16.0	0.0	16.0	17.0	17.0	17.0	15	0.1	0.0
7/1/2021												16	0.0
8/25/2021												18	0.0
9/8/2021	17.0	17.0	18.0	18.0	17.0	0.0	18.0	18.0	18.0	18.0	16	0.3	0.0
10/20/2021												0.0	0.0
11/19/2021												0.0	0.0
12/10/2021	16.0	16.0	17.0	16.0	17.0	0.0	17.0	17.0	17.0	17.0	15	7.6	0.0
1/11/2022											19	0.0	0.0
2/2/2022												0.098	0.0
3/10/2022	15.5	16.5	17.0	16.5	16.5	1.0	16.5	17.0	17.0	17.0	12	0.0	0.0
4/21/2022											19	0.0	0.0
5/16/2022											18	0.0	0.0
6/9/2022	16.0	17.0	17.0	16.0	17.0	0.0	17.0	17.0	17.0	17.0	19	0.0	0.0
7/28/2022											19	1.4	0.0
8/26/2022											19	0.5	0.0
9/22/2022	18.0	18.0	19.0	18.0	18.0	0.0	18.0	19.0	19.0	19.0	18	1.2	0.1
10/13/2022	18.0	18.0	18.0	18.0	18.0	0.0	18.0	18.0	18.0	19.0	19	0.2	0.0
11/7/2022	18.0	18.0	18.0	18.0	18.0	0.0	18.0	18.0	18.0	18.0	19	0.0	0.0
12/9/2022	18.0	18.0	18.0	18.0	18.0	0.0	18.0	18.0	18.0	18.0	19	0.0	0.0
1/31/2023	16.0	17.0	18.0	17.0	17.0	0.0	17.0	18.0	17.0	18.0	18	0.0	0.0
2/21/2023	16.0	17.0	18.0	17.0	17.0	0.0	17.0	18.0	17.0	18.0	18	0.0	0.0
3/10/2023	18.0	18.0	18.0	18.0	18.0	0.0	18.0	18.0	18.0	18.0	19	0.0	0.0
4/6/2023											20	0.0	0.0
5/17/2023											20	0.0	0.0
6/20/2023	17.0	18.0	19.0	18.0	18.0	0.0	18.0	19.0	18.0	19.0	20	0.3	0.1
7/5/2023											20	0.0	0.0
8/17/2023											21	0.0	0.0
9/13/2023	19.0	20.0	20.0	20.0	19.0	0.0	20.0	20.0	20.0	20.0	20	0.0	0.0
10/3/2023											22	0.2	0.3
11/11/2023											20	0.1	0.0
12/12/2023	17.0	18.0	19.0	18.0	19.0	0.0	18.0	20.0	17.0	19.0	20	0.1	0.0
1/12/2024											21	1.4	0.0
2/8/2024											21	1.1	0.0
3/12/2024	17.0	18.0	19.0	18.0	18.0	0.0	18.0	20.0	19.0	19.0	21	0.3	0.0
4/9/2024											22	0.6	0.0
4/15/2024	18.0	19.0	20.0	19.0	18.0	0.2	19.0	20.0	19.0	20.0	23	0.1	0.0
5/8/2024											21	0.1	0.0
6/13/2024	17.0	17.0	18.0	17.0	17.0	0.0	18.0	19.0	17.0	18.0	21	0.0	0.0
7/12/2024											21	0.5	0.2
8/6/2024											21	9.1	5.1
9/6/2024	18.0	19.0	19.0	19.0	18.0	0.0	18.0	20.0	19.0	19.0	21	1.3	0.2

Date	Vapor Monitoring Points (in WC)										
	VMP-1A	VMP-2A	VMP-3A	VMP-4A	VMP-5A	VMP-6A	VMP-7A	VMP-8A	VMP-9A	VMP-9A	
9/26/2019	-0.066	-0.044	-0.075	-0.161	-0.128	+0.000	-0.025	-0.021	N/A	-0.173	
10/3/2019	-0.065	-0.037	-0.053	-0.139	-0.116	+0.000	-0.019	-0.017	N/A	-0.105	
10/9/2019	-0.061	-0.034	-0.045	-0.110	-0.103	+0.000	-0.020	-0.015	N/A	-0.100	
11/5/2019	-0.041	-0.029	-0.023	-0.067	-0.062	+0.010	-0.013	+0.000	N/A	-0.067	
12/3/2019	-0.045	-0.025	-0.031	-0.066	-0.056	+0.020	-0.010	+0.000	N/A	-0.054	
2/11/2020	-0.037	-0.020	-0.015	-0.045	-0.036	+0.015	+0.000	+0.000	N/A	-0.037	
3/27/2020	-0.025	-0.023	-0.016	-0.032	-0.032	+0.010	+0.000	+0.000	N/A	-0.022	
6/28/2020	-0.063	-0.064	-0.063	-0.124	-0.080	NG	-0.010	-0.017	N/A	-0.084	
9/15/2020	-0.063	-0.052	-0.043	-0.093	-0.033	NG	-0.017	-0.014	N/A	-0.058	
12/8/2020	-0.048	-0.033	-0.026	-0.152	-0.05	NG	+0.000	+0.000	N/A	-0.065	
3/30/2021	-0.038	-0.052	-0.032	-0.063	-0.022	NG	-0.020	-0.014	N/A	-0.047	
6/11/2021	-0.073	-0.065	-0.055	-0.105	-0.074	NG	-0.028	-0.022	N/A	-0.074	
9/8/2021	-0.091	-0.088	-0.075	-0.140	-0.086	NG	-0.028	-0.190	N/A	-0.149	
12/10/2021	-0.065	-0.056	-0.043	-0.088	-0.052	NG	-0.017	-0.005	N/A	-0.088	
3/10/2022	-0.045	-0.04	-0.045	-0.080	-0.04	+0.013	-0.010	+0.000	N/A	-0.097	
3/31/2022	NG	NG	NG	NG	NG	NG	NG	+0.000	N/A	NG	
4/21/2022	NG	NG	NG	NG	NG	NG	NG	+0.000	N/A	NG	
5/16/2022	NG	NG	NG	NG	NG	NG	NG	+0.000	N/A	NG	
6/9/2022	-0.068	-0.060	-0.068	-0.097	-0.056	+0.000	-0.027	+0.000	N/A	-0.110	
7/28/2022	NG	NG	NG	NG	NG	NG	NG	-0.018	N/A	NG	
9/22/2022	-0.100	-0.098	-0.105	-0.157	-0.082	+0.000	-0.032	-0.016	N/A	-0.149	
10/13/2022	-0.069	-0.063	-0.071	-0.126	-0.071	+0.000	-0.025	-0.018	N/A	-0.122	
11/7/2022	-0.077	-0.063	-0.084	-0.122	-0.059	+0.000	-0.021	+0.000	N/A	-0.115	
12/9/2022	-0.074	-0.043	-0.046	-0.089	-0.048	+0.000	-0.022	+0.000	N/A	-0.110	
1/31/2023	-0.069	-0.040	-0.042	-0.067	-0.039	+0.000	-0.014	+0.000	N/A	-0.078	
2/21/2023	-0.059	-0.048	-0.061	-0.083	-0.040	+0.000	-0.019	-0.007	N/A	-0.100	
3/10/2023	-0.052	-0.032	-0.054	-0.067	-0.032	+0.000	+0.000	+0.000	N/A	-0.039	
4/12/2023	NG	NG	NG	NG	NG	NG	-0.025	0.000	N/A	NG	
5/17/2023	NG	NG	NG	NG	NG	NG	-0.032	0.000	N/A	NG	
6/20/2023	-0.083	-0.066	-0.085	-0.118	-0.066	+0.000	-0.024	-0.013	N/A	-0.133	
7/5/2023	NG	NG	NG	NG	NG	NG	NG	NG	N/A	NG	
8/17/2023	NG	NG	NG	NG	NG	+0.000	NG	NG	N/A	NG	
9/13/2023	-0.097	-0.079	-0.102	-0.14	-0.083	+0.000	-0.037	-0.013	N/A	-0.140	
10/3/2023	NG	NG	NG	NG	NG	+0.000	NG	NG	N/A	NG	
11/1/2023	NG	NG	NG	NG	NG	+0.000	NG	NG	N/A	NG	
12/12/2023	-0.066	-0.140	-0.203	-0.271	-0.141	+0.000	-0.019	+0.000	N/A	-0.219	
1/12/2024	NG	NG	NG	NG	NG	+0.000	NG	+0.000	N/A	NG	
2/8/2024	NG	NG	NG	NG	NG	+0.000	NG	-0.017	N/A	NG	
3/12/2024	-0.065	-0.045	-0.057	-0.080	-0.039	+0.000	-0.023	+0.000	N/A	-0.084	
4/9/2024	NG	NG	NG	NG	NG	NG	+0.000	-0.030	+0.000	N/A	NG
4/15/2024	-0.077	-0.149	-0.175	-0.092	-0.045	+0.000	-0.027	+0.000	-0.014	-0.274	
5/8/2024	NG	NG	NG	NG	NG	NG	+0.000	NG	+0.000	-0.016	NG
6/13/2024	-0.075	-0.066	Covered	-0.126	-0.061	+0.000	-0.034	+0.000	-0.012	-0.096	
7/1/2024	NG	NG	Covered	NG	NG	NG	NG	-0.014	NG	NG	
8/6/2024	NG	NG	NG	NG	NG	NG	NG	NG	NG	NG	
9/6/2024	-0.116	-0.065	-0.093	-0.154	-0.065	+0.000	-0.059	-0.034	-0.046	-0.130	

**Note:**  
1. Yellow shading indicates that samples did not meet the minimum 0.002 inches WC  
2. Blank space indicates that data was not collected  
3. in WC = inches water column; ppm = parts per million;  
4. N/A = Not Accessible; NG = Not Gauged

**Table 2B**  
**MOD-PAC CORP., 1801 Elmwood Ave, Buffalo, NY**  
**SSDS Post Installation Monitoring Results**  
**Area B - Cold Storage Garage**

Date	Extraction Wells (in WC)								Blower (in WC)	System Effluent PID Reading (ppm)
	EW-1B	EW-2B	EW-3B	EW-4B	EW-5B	EW-6B	EW-7B	EW-8B		
9/26/2019	13.0	13.5	13.5	14.5	13.5	14.0	13.0	12.0	10.5	1.3
10/3/2019	13.0	13.5	13.5	14.0	13.5	14.0	13.0	12.0	10	1.4
10/9/2019	12.5	13.0	13.0	13.5	13.0	13.5	12.0	12.0	10	0.0
11/5/2019	12.0	13.0	12.5	13.0	12.5	13.0	11.5	11.0	9	0.5
12/3/2019	11.0	11.0	11.0	11.5	11.0	11.5	10.5	10.0	8	0.1
1/22/2020										0.0
2/11/2020	12.5	13.0	13.0	13.5	13.0	13.5	12.0	11.5	9	0.0
3/27/2020	14.0	15.0	14.0	15.0	15.0	15.0	14.0	13.5	10	0.0
6/29/2020	16.0	12.0	17.0	12.5	17.0	17.0	16.0	15.5	16	0.0
7/31/2020										0.0
8/28/2020										0.0
9/15/2020	17.0	18.0	17.0	18.0	18.0	18.0	17.0	16.5	16	2.7
10/15/2020										0.3
11/4/2020										0.0
12/8/2020	16.5	17.0	17.0	17.0	17.0	17.0	16.5	16.0	13	0.4
1/4/2021										0.0
2/18/2021										0.0
3/30/2021	16.0	17.0	17.0	17.0	17.0	17.0	16.0	16.0	12	0.0
4/14/2021										0.0
5/20/2021										0.1
6/11/2021	18.0	18.0	19.0	20.0	19.0	19.0	18.0	18.0	18	0.0
7/1/2021									18	0.0
8/25/2021									20	0.0
9/8/2021	20.0	21.0	22.0	23.0	22.0	22.0	21.0	21.0	19	0.0
10/20/2021										0.0
11/19/2021										0.0
12/10/2021	20.0	20.0	21.0	21.0	21.0	21.0	20.0	20.0	16	0.0
1/11/2022									19	0.0
2/2/2022										0.0
3/10/2022	22.0	23.0	23.0	23.5	22.5	23.0	22.5	22.0	20	0.0
4/21/2022									19	0.0
5/16/2022									19	0.0
6/8/2022	26.0	27.0	27.0	28.0	27.0	27.0	27.0	26.0	19	0.0
7/28/2022									25	0.5
8/26/2022									23	0.0
9/22/2022	28.0	29.0	30.0	30.0	29.0	30.0	29.0	28.0	26	2.6
10/13/2022	31.0	32.0	33.0	33.0	32.0	34.0	32.0	32.0	20	0.8
11/7/2022	31.0	32.0	33.0	33.0	33.0	34.0	32.0	32.0	18	0.0
12/8/2022	32.0	33.0	34.0	34.0	33.0	34.0	33.0	32.0	19	0.0
1/31/2023	31.0	32.0	33.0	33.0	32.0	33.0	32.0	32.0	19	0.0
2/21/2023	30.0	31.0	32.0	32.0	31.0	32.0	31.0	30.0	26	0.0
3/10/2023	32.0	32.0	32.0	32.0	32.0	32.0	32.0	32.0	19	0.0
4/6/2023									24	0.0
5/17/2023									29	0.0
6/20/2023	31.0	32.0	32.0	33.0	32.0	33.0	32.0	32.0	30	0.0
7/5/2023									44	0.0
8/17/2023									40	0.0
9/13/2023	37.0	33.0	38.0	36.0	37.0	39.0	37.0	38.0	34	0.0
10/3/2023									34	0.7
11/11/2023									28	0.0
12/12/2023	36.0	37.0	37.0	38.0	37.0	39.0	37.0	37.0	31	0.0
1/12/2024									44	0.2
02/08/2024									45	0.1
3/12/2024	36.0	37.0	37.0	38.0	37.0	39.0	37.0	32.0	31	0.0
4/9/2024	36.0	39.0	37.0	38.0	38.0	39.0		38.0	32	0.0
4/15/2024	36.0	37.0	38.0	38.0	37.0	38.0	37.0	37.0	N/A	0.0
5/8/2024									36	0.1
6/13/2024	37.0	38.0	38.0	39.0	38.0	39.0	38.0	37.0	21	0.0
7/1/2024									24.0	0.0
8/6/2024									38.0	2.7
9/6/2024	39.0	39.0	40.0	40.0	39.0	40.0	39.0	39.0	39.0	0.0

Date	Vapor Monitoring Points (in WC)							
	VMP-1B	VMP-2B	VMP-3B	VMP-4B	VMP-5B	VMP-5BR	VMP-6B	VMP-7B
9/26/2019	N/A	-0.065	-0.419	N/A	-0.044	N/A	-0.016	-0.200
10/3/2019	-0.023	-0.062	-0.303	-0.383	-0.037	N/A	-0.018	-0.196
10/9/2019	-0.018	-0.055	-0.258	-0.329	-0.030	N/A	-0.010	-0.178
11/5/2019	-0.018	-0.018	-0.217	-0.271	-0.014	N/A	+0.000	-0.171
12/3/2019	-0.014	-0.032	-0.114	-0.156	+0.000	N/A	+0.000	-0.136
2/11/2020	+0.000	-0.040	N/A	-0.161	N/A	N/A	+0.000	-0.072
3/27/2020	+0.000	-0.040	-0.163	-0.171	+0.000	N/A	-0.010	-0.152
6/29/2020	-0.018	-0.064	-0.354	-0.343	-0.026	N/A	-0.022	-0.0198
9/15/2020	-0.017	-0.041	-0.118	-0.361	-0.045	N/A	-0.005	-0.160
12/8/2020	+0.000	-0.02	-0.137	-0.208	+0.000	N/A	+0.000	-0.203
3/30/2021	-0.010	-0.045	-0.162	-0.219	+0.000	N/A	-0.010	-0.197
4/14/2021	NG	NG	NG	NG	+0.000	N/A	NG	NG
5/20/2021	NG	NG	NG	NG	-0.014	N/A	NG	NG
6/11/2021	-0.045	-0.051	-0.262	-0.903	-0.039	N/A	-0.016	-0.201
9/8/2021	-0.045	-0.058	-0.285	-1.020	-0.034	N/A	-0.041	-0.060
12/10/2021	-0.010	-0.40	-0.189	-0.177	-0.004	N/A	+0.000	-0.190
1/11/2022	NG	NG	NG	NG	NG	N/A	-0.012	NG
3/10/2022	-0.012	-0.032	-0.141	-0.262	+0.000	N/A	+0.000	-0.133
3/31/2021	NG	NG	NG	NG	-0.167	N/A	-0.014	NG
6/6/2022	-0.014	-0.050	-0.211	-0.299	+0.000	N/A	-0.016	-0.026
7/28/2022	NG	NG	NG	NG	-0.010	N/A	NG	NG
9/22/2022	-0.019	-0.057	-0.238	-0.328	-0.017	N/A	-0.020	-0.263
10/13/2022	-0.045	-0.063	-0.123	-0.215	-0.035	N/A	-0.018	-0.131
11/7/2022	-0.014	-0.057	-0.218	-0.312	+0.000	N/A	-0.016	-0.232
12/8/2022	-0.017	-0.043	-0.153	-0.298	+0.000	N/A	-0.015	-0.156
1/31/2023	-0.009	-0.044	-0.187	-0.279	+0.000	N/A	-0.012	-0.158
2/21/2023	-0.10	-0.045	N/A	-0.299	+0.000	N/A	-0.014	-0.165
3/10/2023	-0.015	-0.030	-0.046	-0.266	+0.000	N/A	-0.015	-0.035
4/12/2023	NG	NG	NG	NG	+0.000	N/A	NG	NG
5/17/2023	NG	NG	NG	NG	+0.000	N/A	NG	NG
6/20/2023	-0.012	-0.045	-0.237	-0.350	+0.000	N/A	-0.017	-0.207
7/5/2023	NG	NG	NG	NG	NG	N/A	NG	NG
8/17/2023	NG	NG	NG	NG	-0.014	N/A	NG	NG
9/13/2023	-0.016	-0.062	-0.433	Covered	-0.011	N/A	-0.018	-0.284
10/3/2023	NG	NG	NG	Covered	NG	N/A	NG	NG
11/11/2023	NG	NG	NG	-0.087	NG	N/A	NG	NG
12/12/2023	-0.016	-0.035	-0.089	-0.319	+0.000	N/A	-0.018	-0.257
1/12/2024	NG	NG	NG	NG	-0.04	N/A	NG	NG
2/8/2024	NG	NG	NG	NG	NG	N/A	NG	NG
3/12/2024	+0.000	-0.001	-0.006	-0.012	+0.000	N/A	+0.000	-0.009
4/9/2024	+0.000	NG	NG	NG	+0.000	N/A	-0.016	NG
4/15/2024	-0.036	-0.101	-0.652	-0.864	+0.000	-0.058	-0.038	-0.695
5/8/2024	-0.012	NG	NG	NG	+0.000	-0.019	NG	NG
6/13/2024	-0.018	-0.047	-0.293	-0.376	+0.000	-0.026	-0.020	-0.290
7/1/2024	NG	NG	NG	NG	-0.010	-0.031	NG	NG
8/6/2024	NG	NG	NG	NG	NG	NG	NG	NG
9/6/2024	-0.018	-0.052	-0.350	-0.373	+0.000	-0.029	-0.015	-0.297

**Note:**

1. Yellow shading indicates that samples did not meet the minimum 0.002 inches WC
2. N/A indicates the VMP was not accessible during the time of the system check
3. Blank space indicates that data was not collected
4. in WC = inches water column; ppm = parts per million;
5. NG = Not Gauged

**Table 2C**  
**MOD-PAC CORP., 1801 Elmwood Ave, Buffalo, NY**  
**SSDS Post Installation Monitoring Results**  
**Area C - Maintenance Area**

Date	Extraction Wells (in WC)			Fan System Effluent PID Reading (ppm)		
	EW-1C	EW-2C	EW-3C	EW-1C	EW-2C	EW-3C
9/26/2019	43.0	40.0		1.4	0.7	
10/3/2019	44.0	45.0		1.0	4.5	
10/9/2019	44.5	45.5		0.0	0.0	
11/5/2019	44.0	46.0		0.0	0.4	
12/3/2019		39.0	28.0		1.2	0.4
1/22/2020					0.4	0.0
2/11/2020	31.0	30.0	27.5	0.2	0.0	0.0
3/27/2020	29.0	32.0	28.0	0.0	0.0	0.0
6/29/2020	27.0	31.0	29.0	0.0	0.0	0.0
7/31/2020				0.0	0.0	0.0
8/28/2020				0.0	0.0	0.0
9/15/2020	28.5	31.0	29.0	0.0	0.0	0.0
10/15/2020				0.0	0.0	0.0
11/4/2020				0.0	0.0	0.0
12/8/2020	31.0	31.0	29.0	0.0	0.0	0.0
1/4/2021				0.0	0.0	0.0
2/18/2021						0.0
3/30/2021		32.0	30.0		0.0	0.0
4/14/2021					0.1	0.0
5/20/2021				0.0	0.0	0.0
6/11/2021	23.0	31.0	30.0	0.0	0.0	0.0
7/1/2021				0.0	0.0	0.0
8/25/2021				0.0	0.0	0.0
9/8/2021	29.0	31.0	30.0	0.0	0.0	0.0
10/20/2021				0.0	0.0	0.0
11/19/2021				0.0	0.0	0.0
12/10/2021	30.0	32.0	30.0	4.7	0.0	0.0
1/11/2022				0.0	0.0	0.0
2/2/2022				0.0	0.0	0.0
3/10/2022	11.0	32.0	31.0	0.0	0.0	0.0
4/21/2022				0.0	0.0	0.0
5/16/2022				0.0	0.0	0.0
6/6/2022	28.0	31.0	32.0	0.0	0.0	0.0
7/28/2022				1.5	0.7	0.1
8/26/2022				0.1	0.0	0.0
9/22/2022	29.0	31.0	32.0	0.0	0.0	0.0
10/13/2022	29.0	31.0	0.0	0.0	0.0	NG
11/7/2022	29.0	31.0	0.0	0.0	0.0	NG
12/9/2022	30.0	30.0	30.0	0.0	0.0	0.0
1/31/2023	0.0	0.0	30.0	NG	NG	0.0
2/21/2023	NG	NG	NG	NG	NG	NG
3/10/2023	0.0	0.0	30.0	0.0	0.0	0.0
4/6/2023	NG	NG	28.0	NG	NG	0.0
5/17/2023	NG	NG	27.0	NG	NG	0.0
6/20/2023	0.0	0.0	29.0	0.0	0.0	0.0
7/5/2023	NG	NG	29.0	NG	NG	0.0
8/17/2023	NG	NG	29.0	NG	NG	0.3
9/13/2023	0.0	0.0	29.0	0.0	0.0	0.0
10/3/2023	35.0	38.0	30.0	27.9	6.7	1.0
11/11/2023	33.0	36.0	29.0	1.1	2.1	0.0
12/12/2023	34.0	37.0	29.0	4.7	2.5	0.1
1/12/2024	34.0	35.0	30.0	2.3	1.8	0.4
2/8/2024	43.0	46.0	30.0	1.6	1.2	0.2
3/12/2024	43.0	46.0	31.0	3.8	2.8	0.5
4/9/2024	44.0	46.0	30.0	0.4	0.0	0.0
4/15/2024	43.0	45.0	30.0	0.0	0.0	0.0
5/8/2024	43.0	45.0	30.0	0.4	0.3	10.3
6/13/2024	42.0	45.0	29.0	1.5	0.8	0.0
7/1/2024	44.0	45.0	30.0	0.1	0.0	0.4
8/6/2024	44.0	45.0	29.0	1.7	1.8	1.1
9/6/2024	42.0	45.0	30.0	0.7	0.8	0.3

Date	Vapor Monitoring Points (in WC)					
	VMP-1C	VMP-2C	VMP-3C	VMP-4C	VMP-10C	VMP-11C
9/26/2019	-0.046	-0.085	+0.000	-0.061		
10/3/2019	-0.055	-0.092	+0.000	-0.081		
10/9/2019	-0.037	-0.075	+0.000	-0.060		
11/5/2019	-0.042	-0.067	+0.000	-0.067		
12/3/2019	+0.000	-0.027	-0.026	+0.004	-0.045	-0.018
2/11/2020	-0.019	-0.026	-0.032	-0.038	-0.045	-0.020
3/27/2020	-0.019	-0.033	-0.038	-0.029	-0.060	-0.021
6/29/2020	-0.019	-0.050	-0.040	-0.018	-0.061	-0.044
9/15/2020	-0.012	-0.040	-0.038	-0.024	-0.039	-0.017
12/8/2020	-0.012	-0.038	-0.026	-0.021	-0.038	-0.016
3/30/2021	+0.000	-0.022	-0.037	+0.000	-0.025	-0.020
6/11/2021	-0.020	-0.054	-0.039	-0.024	-0.058	-0.097
9/8/2021	-0.049	-0.042	-0.040	-0.075	-0.066	-0.022
12/10/2021	-0.026	-0.040	-0.038	-0.021	-0.059	-0.025
2/2/2022	+0.000	-0.028	-0.038	-0.012	-0.034	-0.019
3/10/2022	+0.000	-0.031	-0.038	+0.000	-0.042	-0.022
3/31/2022	-0.021	NG	NG	-0.030	NG	NG
6/6/2022	-0.019	-0.058	-0.037	-0.024	-0.076	-0.039
9/22/2022	-0.021	-0.059	-0.041	-0.018	-0.086	-0.046
10/13/2022	-0.033	-0.042	+0.000	-0.044	-0.044	+0.000
11/7/2022	-0.016	-0.048	+0.000	-0.023	-0.055	+0.000
12/9/2022	-0.041	-0.030	-0.039	-0.045	-0.056	-0.022
1/31/2023	NG	NG	NG	NG	NG	NG
2/21/2023	NG	NG	NG	NG	NG	NG
3/10/2023	+0.000	+0.000	-0.031	+0.000	-0.045	-0.019
4/6/2023	NG	NG	NG	NG	NG	NG
5/17/2023	NG	NG	NG	NG	NG	NG
6/20/2023	+0.000	+0.000	-0.029	+0.000	-0.024	-0.040
7/5/2023	NG	NG	NG	NG	NG	NG
8/17/2023	NG	NG	NG	NG	NG	NG
9/13/2023	+0.000	+0.000	-0.03	+0.000	-0.019	-0.038
10/3/2023	-0.036	-0.063	NG	-0.040	NG	NG
11/11/2023	-0.024	-0.044	-0.046	-0.043	-0.162	-0.108
12/12/2023	-0.016	-0.046	-0.024	-0.028	-0.063	-0.032
1/12/2024	NG	NG	NG	NG	NG	NG
2/8/2024	NG	NG	NG	NG	NG	NG
3/12/2024	-0.051	-0.073	-0.028	-0.069	-0.067	-0.025
4/9/2024	NG	NG	NG	NG	NG	NG
4/15/2024	-0.091	-0.203	-0.059	-0.163	-0.214	-0.078
5/8/2024	NG	NG	NG	NG	NG	NG
6/13/2024	-0.035	-0.081	-0.021	-0.066	-0.102	-0.042
7/1/2024	NG	NG	NG	NG	NG	NG
8/6/2024	NG	NG	NG	NG	NG	NG
9/6/2024	-0.051	-0.094	-0.023	-0.073	-0.124	-0.047

**Note:**

1. Yellow shading indicates that samples did not meet the minimum 0.002 inches WC
2. Blank space indicates that data was not collected
3. in WC = inches water column; ppm = parts per million;
4. N/A = Not Accessible; NG = Not Gauged
5. Please note: The extraction system in area C was operated by fans from 2019-2023. In August 2023, the fans at EW-1C and EW-2C were removed, and a 1.5 hp blower was installed. EW-3C continues to be operated by a fan.



**Table 3**  
**MOD-PAC, Corp. 1801 Elmwood Avenue, Buffalo, NY**  
**Summary of Air Analytical Testing Results**

Parameter	September 2024 - L2453073	
	AREA A-PRE (091624)	AREA A-POST (091624)
<b>Volatile Organic Compounds (ug/m<sup>3</sup>)</b>		
1,1,1-Trichloroethane	ND	ND
1,1,2,2-Tetrachloroethane	ND	ND
1,1,2-Trichloroethane	ND	ND
1,1-Dichloroethane	ND	ND
1,1-Dichloroethene	ND	ND
1,2,4-Trichlorobenzene	ND	ND
1,2,4-Trimethylbenzene	16.4	14.4
1,2-Dibromoethane	ND	ND
1,2-Dichlorobenzene	ND	ND
1,2-Dichloroethane	ND	ND
1,2-Dichloropropane	ND	ND
1,3,5-Trimethylbenzene	4.23	3.6
1,3-Butadiene	ND	ND
1,3-Dichlorobenzene	ND	ND
1,4-Dichlorobenzene	ND	ND
1,4-Dioxane	ND	ND
2,2,4-Trimethylpentane	2.52	1.61
2-Butanone	4.36	ND
2-Hexanone	ND	ND
3-Chloropropene	ND	ND
4-Ethyltoluene	8.7	2.93
4-Methyl-2-pentanone	8.24	6.48
Acetone	195	32.5
Benzene	2.18	3.51
Benzyl chloride	ND	ND
Bromodichloromethane	ND	ND
Bromoform	ND	ND
Bromomethane	ND	ND
Carbon disulfide	3.67	10.3
Carbon tetrachloride	ND	ND
Chlorobenzene	ND	ND
Chloroethane	ND	ND
Chloroform	2.51	ND
Chloromethane	ND	0.968
cis-1,2-Dichloroethene	4.28	ND
cis-1,3-Dichloropropene	ND	ND
Cyclohexane	6.13	1.03
Dibromochloromethane	ND	ND
Dichlorodifluoromethane	2.36	2.32
Ethyl Alcohol	59.7	62.7
Ethyl Acetate	ND	ND
Ethylbenzene	9.69	10.5
Freon-113	ND	ND
Freon-114	ND	ND
Heptane	11.4	2.75
Hexachlorobutadiene	ND	ND
iso-Propyl Alcohol	1120	229
Methyl tert butyl ether	ND	ND
Methylene chloride	ND	3.65
n-Hexane	26	13.2
Naphthalene	ND	ND
o-Xylene	14.4	16.4
p/m-Xylene	43.9	51.3
Styrene	ND	ND
tert-Butyl Alcohol	10.9	ND
Tetrachloroethene	ND	ND
Tetrahydrofuran	ND	ND
Toluene	22	23.7
trans-1,2-Dichloroethene	ND	ND
trans-1,3-Dichloropropene	ND	ND
Trichloroethene	400	ND
Trichlorofluoromethane	5.61	3.64
Vinyl bromide	ND	ND
Vinyl chloride	ND	ND

**Notes:**

- Compounds detected in one or more samples included in this table. For a list of all compounds, refer to analytical report in the Appendix.
- Analytical testing for VOCs via TO-15 completed by Alpha Analytical.
- Results present in ug/m<sup>3</sup> or microgram per cubic meter.
- Parameters shaded in red indicate analytes of concern (Target cVOCs)
- Results in red indicate higher post-carbon readings over pre-carbon readings
- Blank results = No Value Above Detection Limit

Table 4  
MOD-PAC, Corp. 1801 Elmwood Avenue, Buffalo, NY  
Summary of Air Analytical Testing Results

Parameter	October 2019 - L1946093			November 2019 - L1952487			December 2019 - L1957660			February 2020 - L2006152			June 2020 - L2027736		September 2020 - L2038512		09/23/2020	December 2020 - L2054640		March 2021 - L2115934		June 2021 - L2131935		September 2021 - L2148116		12/10/2021	December 2021 - L2168195		March 2022 - L2212728		
	AREA A- PRE	AREA A- POST	AREA B	AREA A- PRE (110519)	AREA A- POST (110519)	AREA B (110519)	AREA A- PRE (120319)	AREA A- POST (120319)	AREA B (120319)	AREA A- PRE (021120)	AREA A- POST (021120)	AREA B (120319)	AREA A- PRE (063020)	AREA A- POST (063020)	AREA A- PRE (091520)	AREA A- POST (091520)		AREA A- PRE (120820)	AREA A- POST (120820)	AREA A- PRE (033021)	AREA A- POST (033021)	AREA A- PRE (061121)	AREA A- POST (061121)	AREA A- PRE (090821)	AREA A- POST (090821)		AREA A- PRE (121021)	AREA A- POST (121021)	AREA A- PRE (031022)	AREA A- POST (031022)	
<b>Volatile Organics in Air (ug/m<sup>3</sup>)</b>																															
1,1,1-Trichloroethane	1.11	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,1,2,2-Tetrachloroethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,1,2-Trichloroethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,1-Dichloroethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,1-Dichloroethane	94.8	ND	4.52	35.5	ND	ND	41.6	5.55	0.979	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2,4-Trichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2,4-Trimethylbenzene	2.5	ND	ND	ND	ND	ND	ND	ND	ND	48.5	30.2	56	21.8	21.5	64.4	63.4		29.7	23.7	34.4	28.8	46.1	38.9	42.4	53.1	59	49.2	7.28	4.56		
1,2-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,2-Dichloropropane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,3,5-Trimethylbenzene	1	ND	ND	ND	ND	ND	ND	ND	ND	7.87	4.7	10.2	5.7	4.75	14.5	17.2		8.95	6.44	12.4	9.54	14.2	11.2	10.2	13.6	21.3	17.2	2.36	1.43		
1,3-Butadiene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,3-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,4-Dichlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
1,4-Dioxane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
2,2,4-Trimethylpentane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.976	2.98	ND	ND		ND	ND	ND	ND	3.14	ND	ND	1.37	1.37	ND	ND	ND	ND	
2-Butanone	9.88	ND	3.07	4.13	ND	ND	5.28	ND	ND	4.04	ND	ND	6.25	2.45	ND	ND		2.16	ND	2.98	ND	3.89	ND	2.53	ND	2.78	1.68	1.8	ND	ND	
2-Hexanone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
3-Chloropropene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
4-Ethyltoluene	ND	ND	ND	ND	ND	ND	ND	ND	ND	14.5	9.49	21.8	4.22	3.87	12.4	10.9		3.95	2.79	6.1	4.46	10.7	8.26	6	8.26	30	21.6	ND	ND		
4-Methyl-2-pentanone	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	9.71	ND	4.47	ND	ND	3.53	ND	ND	ND	ND	ND	
Acetone	50.2	10.5	22.7	49.9	ND	ND	69.8	75.5	4.44	13.3	87.4	ND	53.4	100	10.6	26.5		195	12.3	73.6	12.5	73.6	20.7	38.2	40.4	108	22.2	134	ND	ND	
Benzene	0.891	ND	ND	ND	ND	ND	ND	ND	ND	5.34	2.5	10.4	ND	0.987	4.79	2.43		1.42	0.69	2.25	1.03	10.7	4.98	2.75	5.46	2.58	1.04	ND	ND	ND	
Benzyl chloride	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Bromodichloromethane	ND	ND	ND	9.71	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Bromoforn	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Bromomethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Carbon disulfide	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.935	ND	ND	21.5	ND	5.82	6.42		1.45	0.931	2.42	0.944	7.41	2.68	3.83	12.5	4.61	2.56	1.3	0.956		
Carbon tetrachloride	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.26	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Chlorobenzene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Chloroethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Chloroform	14.4	ND	ND	9.86	ND	ND	20.3	1.69	ND	17	1.51	ND	16.7	31.8	20.7	17.5		27.1	1.35	38.4	12.6	46.7	59.6	31.5	42.7	26.2	1.2	40.5	0.96		
Chloromethane	0.591	0.745	ND	ND	ND	ND	ND	0.603	0.785	ND	0.446	1.21	ND	0.77	ND	0.438		0.626	0.630	0.648	0.766	ND	0.558	ND	0.564	0.605	0.465	0.62	1.01		
cis-1,2-Dichloroethene	88.8	ND	ND	33.5	ND	ND	41.6	5.55	0.979	22.5	12.5	ND	26.1	63	19.2	21.7		15.1	ND	11.2	11.3	11.7	29.1	10.1	13.7	3.87	ND	3.26	ND	ND	
cis-1,3-Dichloropropene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Cyclohexane	4.23	ND	2	ND	2.52	ND	ND	ND	ND	1.61	ND	0.847	ND	ND	2.54	0.823		2.1	ND	1.41	ND	2.42	ND	ND	1.29	1.61	ND	ND	ND	ND	
Dibromochloromethane	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Dichlorodifluoromethane	1.99	1.78	1.98	2.13	ND	ND	ND	2.1	2.93	ND	1.47	1.99	ND	2.15	ND	1.61		2.41	2.38	1.95	2.04	2.06	1.87	2.64	2.14	2.1	ND	2.35	2.39		
Ethyl Alcohol	14.3	23.4	16	22.2	ND	61.6	43.5	34.5	10.3	63.7	40.9	30.1	143	112	106	81.8		91	57.1	71.6	86.7	87.8	61.6	49.7	64.1	79	23.2	129	ND	ND	
Ethyl Acetate	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	3.27	3.13	4.4	4.14	ND	ND	3.41	2.5	ND	ND	ND	
Ethylbenzene	1.58	ND	0.973	2.32	ND	ND	3.54	ND	ND	37.6	20	60.4	6.65	5.13	17.9	13.6		16.8	5.08	15.9	6.91	19.1	11.5	9.64	16.8	7.12	4.17	3.61	ND	ND	
Freon-113	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Freon-114	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	
Heptane	14.3	ND	2.35	9.51	ND	6.27	18.2	ND	1.25	16.6	1.01	14.1	5.7	1.25	6.31	1.31		24.9	ND	7.38	0.836	6.64	1.94	1.98	3.74	7.09	ND	13.2	ND	ND	
Hexachlorobutadiene	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
iso-Propyl Alcohol	44	48.2	28	103	ND	742	275	1.96	7.03	157	9.44	44.2	191	472	83.8	34.4		371	32.9	253	164	95.9	533	38.8	95.9	256	16.1	283	3.22	ND	
Methyl tert butyl ether	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND
Methylene chloride	9.21	13.2	9.87	3.68	5.45	5.35	ND	4.45	3.61	ND	ND	ND	ND	ND	ND	ND		ND	ND	ND	1.79	ND	ND	6.62	ND	ND	ND	1.75	ND	ND	ND
Naphthalene	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA		NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA
n-Heptane	6.06	5.08	1.72	5.22	1.89	3.98	28.2	1.2	1.54	20.7	0.948	6.1	12.2	2.59	29.3	3.67		18.1	2.31	33.7	5.15	73.7	14.9	4.12							

Table 4  
MOD-PAC, Corp. 1801 Elmwood Avenue, Buffalo, NY  
Summary of Air Analytical Testing Results

Parameter	June 2022- L2229574		September 2022 L2252350		12/09/2022	December 2022 - L2269445		March 2023 - L2312615		June 2023 - L2335506		September 2023 - L2353358		December 2023 - L2373355		March 2024 - L2413550		June 2024 - L2434419		07/18/2024	September 2024 - L2453073	
	AREA A- PRE (060622)	AREA A- POST (060622)	AREA A- PRE (092222)	AREA A- POST (092222)		AREA A- PRE (120922)	AREA A- POST (120922)	AREA A- PRE (030823)	AREA A- POST (030823)	AREA A- PRE (062023)	AREA A- POST (062023)	AREA A- PRE (091323)	AREA A- POST (091323)	AREA A- PRE (121223)	AREA A- POST (121223)	AREA A- PRE (031224)	AREA A- POST (031224)	AREA A- PRE (061824)	AREA A- POST (061824)		AREA A- PRE (091624)	AREA A- POST (091624)
Volatile Organics in Air (ug/m <sup>3</sup> )																						
1,1,1-Trichloroethane	ND	ND	ND	ND	CARBON CHANGEOUT	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	CARBON CHANGEOUT	ND	ND
1,1,2,2-Tetrachloroethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,1,2-Trichloroethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,1-Dichloroethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,1-Dichloroethene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,2,4-Trichlorobenzene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,2,4-Trimethylbenzene	ND	9.83	4.33	4.39		2.89	3.58	2.16	ND	5.8	4.78	4.35	3.34	4.24	2.3	10.9	3.06	9.98	23.2		16.4	14.4
1,2-Dibromochloroethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,2-Dichlorobenzene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,2-Dichloroethane	ND	0.999	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,2-Dichloropropane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,3,5-Trimethylbenzene	ND	2.7	1.33	1.23		ND	1.55	ND	ND	1.7	1.24	1.15	ND	1.59	ND	4.48	1.26	2.79	5.9		4.23	3.6
1,3-Butadiene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,3-Dichlorobenzene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,4-Dichlorobenzene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
1,4-Dioxane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
2,2,4-Trimethylpentane	ND	ND	1.22	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.31	ND	ND	ND		2.52	1.61
2-Butanone	ND	3.27	2.92	3.16		2.08	ND	4.13	ND	4.98	1.79	4.16	2.01	1.86	ND	2.66	ND	6.64	ND		4.36	ND
2-Hexanone	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
3-Chloropropene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
4-Ethyltoluene	ND	1.85	ND	ND		ND	ND	ND	ND	1.23	ND	ND	ND	0.998	ND	2.82	ND	2.19	6.19		8.7	2.93
4-Methyl-2-pentanone	ND	ND	ND	3.43		ND	ND	ND	ND	2.42	ND	5.49	ND	ND	ND	ND	ND	15.9	24.8		8.24	6.48
Aceetone	688	58.7	35.6	35.5		196	17.3	466	23.6	112	19.1	62.5	15.9	80.8	14.3	236	37.9	143	130		195	22.5
Benzene	ND	1.53	1.56	ND		1.83	0.757	1.45	ND	1.8	1.04	1.29	0.652	0.706	ND	1.04	ND	1.87	5.81		2.18	3.51
Benzyl chloride	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Bromodichloromethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Bromochloromethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Bromomethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Carbon disulfide	7.51	3.74	8.16	6.26		4.20	0.782	ND	3.21	7.29	2.3	6.17	1.89	3.89	1.4	0.772	0.866	5.89	24.2		3.67	10.3
Carbon tetrachloride	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Chlorobenzene	ND	ND	0.953	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Chloroethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Chloroform	21.6	1.67	14	31.3		24.4	ND	18.9	ND	15.1	ND	3.97	15.1	3.63	9.96	1.87	4.44	5.57	26.2		2.51	ND
Chloromethane	ND	0.812	0.849	0.518		0.748	0.791	ND	ND	0.772	0.776	0.653	0.586	0.69	0.578	1.39	0.64	2.03	3.63		ND	0.968
cis-1,2-Dichloroethene	ND	0.999	5.27	6.03		3.30	ND	3.71	ND	5.15	1.34	5.19	4.32	3.85	3.16	2.34	4.24	6.54	24.8		4.28	ND
cis-1,3-Dichloropropene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Cyclohexane	ND	ND	0.981	ND		0.898	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		6.13	1.03
Dibromochloromethane	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Dichlorodifluoromethane	ND	3.12	3.2	2.27		2.61	ND	2.53	2.84	3.19	2.91	2.33	1.84	2.62	2.2	2.19	2.88	4.03	7.22		2.36	2.32
Ethyl Alcohol	148	119	126	83.8		127	25.1	114	121	61	57.8	51.8	43	28.6	52	107	87.1	50.9	166		59.7	62.7
Ethyl Acetate	ND	3.6	4.72	ND		170	137	214	170	178	176	50.1	47.2	24.2	38.9	46.1	41.4	ND	ND		ND	ND
Ethylbenzene	ND	3.87	2.21	1.12		3.86	1.21	2.68	ND	4.08	2.24	2.42	1.23	1.52	ND	3.16	1.55	4.2	13.2		9.69	10.5
Freon-113	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Freon-114	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Heptane	ND	1.75	1.79	ND		9.02	ND	18	ND	1.36	ND	1.23	ND	2.19	ND	8.11	0.91	3.75	4.67		11.4	2.75
Hexachlorobutadiene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
iso-Propyl Alcohol	5090	733 D	56.5	157		467	50.9	637	280	213	551	94.9	317	96.8	160	217	438	178	1150		1120	229
Methyl tert butyl ether	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Methylene chloride	ND	ND	3.07	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	3.65
Naphthalene	NA	NA	NA	NA		NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	NA	ND	4.72	18.6		26	13.2
n-Hexane	14.4	4.86	12.5	8.07		27.7	4.44	ND	ND	9.8	7.08	13.5	10.3	6.8	1.74	11.1	ND	ND	ND		ND	ND
p-Xylene	ND	6.34	3.61	2.28		4.60	2.33	3.03	ND	5.73	4.05	3.76	2.22	2.12	1.32	3.61	1.76	6.39	19.7		14.4	16.4
p/m-Xylene	18.6	17.3	9.86	5.26		14.8	6.30	10.6	ND	18.2	11.60	11	6.08	6.47	3.69	11.9	5.95	19.3	63.9		43.9	51.3
Styrene	ND	0.856	ND	ND		1.26	ND	ND	ND	1.91	0.975	1.32	ND	ND	ND	ND	2.27	ND	ND		ND	ND
tert-Butyl Alcohol	20.3	ND	6.55	4.79		16.6	ND	18	ND	4.18	ND	4.55	2.37	3.3	4.37	8.61	11.1	20.6	49.1		10.9	ND
Tetrachloroethene	ND	ND	2.31	ND		2.94	5.51	4.17	ND	2.27	ND	1.67	ND	ND	ND	ND	9.97	27.7	ND		ND	ND
Tetrahydrofuran	ND	4.16	ND	2.22		ND	ND	ND	ND	2.14	ND	1.91	ND	ND	ND	ND	2.48	5.75	ND		ND	ND
Toluene	20.3	18.4	11.6	4.37		18.6	4.33	10.9	3.66	15.2	8.89	8.48	5.2	4.52	2.86	11	4.6	17.6	47.1		22	23.7
trans-1,2-Dichloroethene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	3.1	ND		ND	ND
trans-1,3-Dichloropropene	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Trichloroethene	262	18.4	353	29.4		250	8.38	183	ND	327	18.3	313	18.4	248	12.6	152	9.24	450	20		400	ND
Trichlorofluoromethane	ND	5.22	3.73	4.61		1.48	ND	ND	ND	4.81	7.31	5.68	4.78	2.16	1.73	2.81	2.12	10.2	25		5.61	3.64
Vinyl bromide	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Vinyl chloride	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	ND		ND	ND
Total Target cVOCs	262.00	19.40	363.65	35.43		256.24	13.89	190.88	0.00	334.42	19.64	319.86	22.72	251.85	15.76	154.34	13.48	466.51	72.50		404.28	3.65
Percent Decrease of																						

Table 5  
Historical Groundwater Monitoring Data Summary  
MOD-PAC CORP.

Monitoring Well	Date	Top of Casing (ft)	Depth to Water (ft)	GW Elevation (ft)	Trichloroethene (µg/L) NY-TGGS-GA (5 µg/L)	% Increase/Decrease TCE
MW - 3	2/5/18	600.71	5.05	595.66	280	Baseline
	Potassium Permanganate Pilot Study June 27, 2019 - June 28, 2019					
	7/16/19	600.71	NG	NG	ND	-100.00
	Potassium Permanganate Injections October 1, 2019 - October 10, 2019					
	10/24/19	600.71	NG	NG	220	-21.43
	4/15/20	600.71	5.54	595.17	370 JH	32.14
	3/10/21	600.71	5.10	594.61	NT	N/A
	3/30/21	600.71	5.35	594.76	NT	N/A
	4/14/21	600.71	5.98	594.73	340	21.43
	5/20/21	600.71	5.10	594.61	NT	N/A
	6/11/21	600.71	6.12	594.59	NT	N/A
	7/1/21	600.71	6.30	594.41	400	42.86
	8/25/21	600.71	5.80	594.91	NT	N/A
	9/22/21	600.71	5.45	595.26	NT	N/A
	11/19/21	600.71	5.30	595.41	340	21.43
	12/10/21	600.71	5.55	595.16	NT	N/A
	1/12/22	600.71	5.70	595.01	190	-32.14
	2/2/22	600.71	6.09	594.62	NT	N/A
	3/10/22	600.71	6.44	594.27	NT	N/A
	4/5/22	600.71	5.65	595.06	280	0.00
	5/16/22	600.71	5.81	594.90	NT	N/A
	6/6/22	600.71	5.70	595.01	NT	N/A
	7/6/22	600.71	5.91	594.80	240	-14.29
	8/9/22	600.71	5.85	594.86	NT	N/A
	9/22/22	600.71	6.18	594.53	NT	N/A
	10/7/22	600.71	6.03	594.68	350	25.00
	11/7/22	600.71	5.71	595.00	NT	N/A
	12/8/22	600.71	5.55	595.15	NT	N/A
	1/5/23	600.71	4.70	596.01	170	-39.29
	2/21/23	600.71	5.70	595.01	NT	N/A
	3/24/23	600.71	5.41	595.30	NT	N/A
	4/6/23	600.71	5.35	595.36	120 J	-57.14
	5/17/23	600.71	5.80	594.91	NT	N/A
	6/20/23	600.71	7.18	593.53	NT	N/A
	7/25/23	600.71	NG	NG	NT	N/A
	8/17/23	600.71	5.95	594.76	NT	N/A
	10/3/23	600.71	6.30	594.41	400	42.86
	1/12/24	600.71	5.28	595.43	330	17.86
	4/9/24	600.71	5.62	595.09	390	7.14
	7/1/24	600.71	5.26	595.45	250	-10.71
MW - 11	2/5/18	600.41	4.66	595.75	40	Baseline
	Potassium Permanganate Pilot Study June 27, 2019 - June 28, 2019					
	7/16/19	600.41	NG	NG	ND	-50.00
	Potassium Permanganate Injections October 1, 2019 - October 10, 2019					
	10/24/19	600.41	NG	NG	16	-60.00
	4/15/20	600.41	5.21	595.14	45 JH	12.50
	3/10/21	600.41	5.82	594.59	NT	N/A
	3/30/21	600.41	5.74	594.67	NT	N/A
	4/14/21	600.41	5.74	594.67	16	-60.00
	5/20/21	600.41	5.84	594.57	NT	N/A
	6/11/21	600.41	5.85	594.56	NT	N/A
	7/1/21	600.41	6.00	594.41	47	17.50
	8/25/21	600.41	5.58	594.83	NT	N/A
	9/22/21	600.41	5.32	595.09	NT	N/A
	11/19/21	600.41	5.15	595.26	32	-20.00
	12/10/21	600.41	5.35	595.06	22	N/A
	1/12/22	600.41	5.45	594.95	22	-45.00
	2/2/22	600.41	5.80	594.61	NT	N/A
	3/10/22	600.41	5.21	595.20	NT	N/A
	4/5/22	600.41	5.41	594.96	24	-40.00
	5/16/22	600.41	5.49	594.92	NT	N/A
	6/6/22	600.41	5.46	594.95	NT	N/A
	7/6/22	600.41	5.63	594.78	27	-32.50
	8/9/22	600.41	5.71	594.70	NT	N/A
	9/22/22	600.41	5.90	594.51	NT	N/A
	10/7/22	600.41	5.80	594.61	34	-15.00
	11/7/22	600.41	5.61	594.80	NT	N/A
	12/8/22	600.41	5.38	595.03	31	N/A
	1/5/23	600.41	4.73	595.68	31	-22.50
	2/21/23	600.41	5.50	594.91	NT	N/A
	3/24/23	600.41	5.39	595.02	NT	N/A
	4/6/23	600.41	4.60	595.81	19	-52.50
	5/17/23	600.41	5.60	594.81	NT	N/A
	6/20/23	600.41	5.94	594.47	NT	N/A
	7/25/23	600.41	5.60	594.81	23	-42.50
	8/17/23	600.41	5.74	594.67	NT	N/A
	10/3/23	600.41	6.05	594.36	12	-10.00
	1/12/24	600.41	5.34	595.07	12	-70.00
	4/9/24	600.41	5.58	594.83	29	-27.50
	7/1/24	600.41	5.30	595.11	37	7.50
MW - 12	2/5/18	600.50	4.52	595.98	0.44 J	Baseline
	Potassium Permanganate Pilot Study June 27, 2019 - June 28, 2019					
	7/16/19	600.50	NG	NG	ND	-100.00
	Potassium Permanganate Injections October 1, 2019 - October 10, 2019					
	10/24/19	600.50	NG	NG	ND	-100.00
	4/15/20	600.50	4.41	596.09	ND	-100.00
	3/10/21	600.50	5.03	595.47	NT	N/A
	3/30/21	600.50	4.86	595.64	NT	N/A
	4/14/21	600.50	4.86	595.64	ND	-100.00
	5/20/21	600.50	5.05	595.45	NT	N/A
	6/11/21	600.50	5.10	595.40	NT	N/A
	7/1/21	600.50	5.35	595.15	ND	-100.00
	8/25/21	600.50	4.80	595.70	NT	N/A
	9/22/21	600.50	4.40	596.10	NT	N/A
	11/19/21	600.50	4.10	596.40	ND	-100.00
	12/10/21	600.50	4.35	596.15	NT	N/A
	1/12/22	600.50	4.58	595.92	ND	-100.00
	2/2/22	600.50	5.20	595.30	NT	N/A
	3/10/22	600.50	4.30	596.20	NT	N/A
	4/5/22	600.50	4.41	596.09	ND	-100.00
	5/16/22	600.50	5.30	595.20	NT	N/A
	6/6/22	600.50	4.73	595.77	NT	N/A
	7/6/22	600.50	4.10	596.40	ND	-100.00
	8/9/22	600.50	4.89	595.61	NT	N/A
	9/22/22	600.50	5.15	595.35	NT	N/A
	10/7/22	600.50	5.04	595.46	ND	-100.00
	11/7/22	600.50	4.62	595.88	NT	N/A
	12/8/22	600.50	3.54	596.96	NT	N/A
	1/5/23	600.50	3.42	596.96	ND	-100.00
	2/21/23	600.50	4.55	595.95	NT	N/A
	3/24/23	600.50	3.38	596.11	NT	N/A
	4/6/23	600.50	3.78	595.74	ND	-100.00
	5/17/23	600.50	4.69	595.81	NT	N/A
	6/20/23	600.50	5.20	595.30	NT	N/A
	7/25/23	600.50	4.71	595.79	0.20 J	-54.55
	8/17/23	600.50	4.94	595.56	NT	N/A
	10/3/23	600.50	5.39	595.11	0.18 J	-59.09
	1/12/24	600.50	4.14	596.36	NT	N/A
	4/9/24	600.50	4.41	596.09	ND	N/A
	7/1/24	600.50	4.10	596.40	0.86	85.45
MW - 13	2/5/18	600.31	4.44	595.87	160	Baseline
	Potassium Permanganate Pilot Study June 27, 2019 - June 28, 2019					
	7/16/19	600.31	NG	NG	ND	-51.25
	Potassium Permanganate Injections October 1, 2019 - October 10, 2019					
	10/24/19	600.31	NG	NG	240	50.00
	4/15/20	600.31	3.70	596.61	140 JH	-12.50
	3/10/21	600.31	4.25	596.06	NT	N/A
	3/30/21	600.31	4.10	596.21	NT	N/A
	4/14/21	600.31	4.13	596.19	95	-40.03
	5/20/21	600.31	4.32	595.99	NT	N/A
	6/11/21	600.31	4.40	595.91	NT	N/A
	7/1/21	600.31	4.60	595.71	150	-6.25
	8/25/21	600.31	4.10	596.21	NT	N/A
	9/22/21	600.31	3.35	596.95	NT	N/A
	11/19/21	600.31	3.30	597.01	73	-54.38
	12/10/21	600.31	3.50	596.81	NT	N/A
	1/12/22	600.31	3.85	596.46	74	-53.75
	2/2/22	600.31	4.30	596.01	NT	N/A
	3/10/22	600.31	4.46	595.85	NT	N/A
	4/5/22	600.31	3.80	596.51	89	-63.13
	5/16/22	600.31	4.10	596.21	NT	N/A
	6/6/22	600.31	4.23	596.08	NT	N/A
	7/6/22	600.31	4.11	596.20	89	-44.38
	8/9/22	600.31	3.90	596.41	NT	N/A
	9/22/22	600.31	4.45	595.86	NT	N/A
	10/7/22	600.31	5.66	594.65	72	-55.00
	11/7/22	600.31	3.78	596.53	NT	N/A
	12/8/22	600.31	3.45	596.86	NT	N/A
	1/5/23	600.31	2.62	597.69	35	-78.13
	2/21/23	600.31	3.81	596.50	NT	N/A
	3/24/23	600.31	3.46	596.85	NT	N/A
	4/6/23	600.31	3.10	597.21	32 J	-80.00
	5/17/23	600.31	4.01	596.30	NT	N/A
	6/20/23	600.31	5.50	594.61	NT	N/A
	7/25/23	600.31	3.98	596.33	90	-43.75
	8/17/23	600.31	4.20	596.11	NT	N/A
	10/3/23	600.31	6.70	593.61	71	-55.03
	1/12/24	600.31	3.11	597.20	36	-77.50
	4/9/24	600.31	3.65	596.65	43	-73.13
	7/1/24	600.31	3.27	597.04	44	-66.42
MW - 14	3/10/21		6.76	-6.76	NT	N/A
	3/30/21		6.72	-6.72	NT	N/A
	4/14/21		6.73	-6.73	NT	N/A
	5/20/21		6.75	-6.75	NT	N/A
	6/11/21		6.80	-6.80	NT	N/A
	7/1/21		6.95	-6.95	NT	N/A
	8/25/21		6.50	-6.50	NT	N/A
	9/22/21		6.15	-6.15	NT	N/A
	11/19/21		6.10	-6.10	NT	N/A
	12/10/21		6.30	-6.30	NT	N/A
	1/12/22		6.40	-6.40	NT	N/A
	2/2/22		6.74	-6.74	NT	N/A
	3/10/22		7.36	-7.36	NT	N/A
	4/5/22		6.40	-6.40	NT	N/A
	5/16/22		6.54	-6.54	NT	N/A
	6/6/22		6.31	-6.31	NT	N/A
	7/6/22		6.5			

Table 6  
Historical Groundwater Monitoring and Sampling Data Summary  
MOD-PAC CORP.

Monitoring Well	Date	Top of Casing (ft)	Depth to Water (ft)	GW Elevation (ft)	1,1-Dichloroethene (µg/L)	2-Butanone (µg/L)	Acetone (µg/L)	Benzene (µg/L)	cis-1,2-Dichloroethene (µg/L)	trans-1,2-Dichloroethene (µg/L)	Trichloroethene (µg/L)	Vinyl chloride (µg/L)	Total VOCs (µg/L)	% Increase/Decrease TCE
MW - 3	2/5/18	600.71	5.05	595.66	ND	ND	ND	ND	89	14	280	13	387.0	Baseline
	7/16/19	600.71	NG	NG	ND	3.10 J	38	ND	ND	ND	ND	ND	43.4	-100.00
	10/24/2019 <sup>1</sup>	600.71	NG	NG	ND	ND	<20	<1	30	3	220	<1	253.0	-21.43
	4/15/20	600.71	5.54	595.17	ND	ND	6.40 J	ND	57	7.3	370 JH	3.7	444.4	32.14
	4/14/21	600.71	5.88	594.73	0.88 J	ND	ND	ND	82	8.8	340	5.6	440.5	21.43
	7/1/21	600.71	6.30	594.41	2.0	ND	ND	0.41 J	140	16	400	8.1	566.5	42.86
	11/19/21	600.71	5.30	595.41	0.77 J	ND	ND	ND	43	4 J	340	2.9	390.7	21.43
	1/12/22	600.71	5.70	595.01	0.86	ND	ND	0.16 J	57	3.3	190	3.5	254.8	-32.14
	4/5/22	600.71	5.65	595.06	0.44 J	ND	ND	ND	46	5.1 J	280	2.3 J	333.8	0.00
	7/6/22	600.71	5.91	594.80	0.48 J	ND	ND	ND	74	6.2	240	3.7	324.4	-14.29
	10/7/22	600.71	6.03	594.69	0.76 J	6.50 J	7.60 J	0.34 J	92	6.5	350	7.2	470.9	25.00
	1/5/23	600.71	4.70	596.01	0.24 J	ND	ND	ND	29	1.5 J	170 R1	0.55 J	201.3	-39.29
	4/6/23	600.71	5.35	595.36	ND	ND	ND	ND	17 J	0.92 J	120 J	0.41 J	138.3	-57.14
	7/25/23	600.71	NG	NG	NT	NT	NT	NT	NT	NT	NT	NT	NT	NT
	10/3/23	600.71	6.30	594.41	ND	ND	ND	ND	99	8.3 J	400	4.8	512.1	42.86
	1/12/24	600.71	5.28	595.43	0.35 J	ND	ND	ND	66	5.4	330	1.7 J	403.5	17.86
	4/9/24	600.71	5.62	595.09	0.41 J	ND	ND	ND	54	4.9 J	300	1.9 J	361.2	7.14
	7/1/24	600.71	5.26	595.45	ND	ND	ND	ND	66	5.2 J	250	1.9 J	323.1	-10.71
	2/5/18	600.41	4.66	595.75	ND	2.3 J	9.4	0.16 J	3.1	2.9	40	5.6	64.56	Baseline
MW - 11	7/16/19	600.41	NG	NG	0.35 J	4.15	ND	ND	ND	ND	20	9.8	73.65	-50.00
	10/24/2019 <sup>1</sup>	600.41	NG	NG	ND	150 J	920	<10	<10	<10	16	ND	1086.0	-60.00
	4/15/20	600.41	5.27	595.14	ND	2.2 J	11	0.21 J	7	10	45 JH	9	84.4	12.50
	4/14/21	600.41	5.74	594.67	ND	ND	ND	ND	8	9.4	16	5.7	39.1	-60.00
	7/1/21	600.41	6.00	594.41	0.35 J	ND	ND	0.25 J	13	10	87.0	10	87.0	17.50
	11/19/21	600.41	5.15	595.26	0.47 J	ND	ND	0.25 J	17	30	32	7.8	87.3	-20.00
	1/12/22	600.41	5.45	594.96	0.31 J	ND	ND	0.20 J	11	19	22	6.2	58.7	-45.00
	4/5/22	600.41	5.45	594.96	0.27 J	ND	ND	0.17 J	9.8	15	24	9.7	58.9	-40.00
	7/6/22	600.41	5.63	594.78	0.36 J	ND	3.6 J	0.22 J	15	20	27	10	76.2	-32.50
	10/7/22	600.41	5.80	594.61	ND	ND	ND	0.22 J	15	34	7.2	34	69.4	-15.00
	1/5/23	600.41	4.73	595.68	0.25 J	ND	ND	0.16 J	11	16	31	9.4	67.8	-22.50
	4/6/23	600.41	4.60	595.81	0.39 J	ND	ND	ND	10 J	16	19 J	10	55.4	-52.50
	7/25/23	600.41	5.60	594.81	0.22 J	ND	2.5 J	0.2 J	12	17	23	17	71.9	-42.50
	10/3/23	600.41	6.05	594.36	ND	ND	5.7	ND	11	12	12	8.5	49.2	-70.00
	1/12/24	600.41	5.34	595.07	0.22 J	ND	ND	ND	13	12	8.7	44.9	70.00	-70.00
	4/9/24	600.41	5.58	594.83	0.52	ND	2.4 J	0.17 J	12	18	25	12	74.1	-27.50
	7/1/24	600.41	5.30	595.11	0.22 J	ND	3.1 J	0.23 J	17	21	37	14	92.6	-7.50
MW - 12	2/5/18	600.50	4.52	595.98	ND	ND	2.2 J	ND	ND	ND	0.44 J	ND g	2.64	Baseline
	7/16/19	600.50	NG	NG	ND	ND	3 J	ND	ND	ND	ND	ND	3.0	-100.00
	10/24/2019 <sup>1</sup>	600.50	NG	NG	ND	ND	<200	ND	ND	ND	ND	ND	ND	-100.00
	4/15/20	600.50	4.41	596.09	ND	ND	11	ND	ND	ND	ND	ND	11.0	-100.00
	4/14/21	600.50	4.86	595.64	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	7/1/21	600.50	5.35	595.15	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	11/19/21	600.50	4.10	596.40	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	1/12/22	600.50	4.58	595.92	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	4/5/22	600.50	4.41	596.09	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	7/6/22	600.50	4.10	596.40	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	10/7/22	600.50	5.04	595.46	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	1/5/23	600.50	3.54	596.96	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	4/6/23	600.50	3.76	596.74	ND	ND	ND	ND	ND	ND	ND	ND	ND	-100.00
	7/25/23	600.50	4.71	595.79	ND	ND	ND	ND	ND	ND	0.20 J	ND	0.2	-54.55
	10/3/23	600.50	5.39	595.11	ND	ND	ND	ND	ND	ND	0.18 J	ND	0.2	-59.09
	1/12/24	600.50	4.14	596.36	ND	ND	ND	ND	ND	ND	ND	ND	0.0	-100.00
	4/9/24	600.50	4.41	596.09	ND	ND	ND	ND	ND	ND	ND	ND	0.0	-100.00
	7/1/24	600.50	4.10	596.40	ND	ND	ND	ND	ND	ND	0.86	ND	0.9	95.45
MW - 13	2/5/18	600.31	4.44	595.87	1	ND	ND	ND	180	4.1	160	25	371.3	Baseline
	7/16/19	600.31	NG	NG	1.20 J	ND	ND	ND	400	3.9 J	78	56	539.1	-51.25
	10/24/2019 <sup>1</sup>	600.31	NG	NG	<1	ND	28	ND	97	2	240	2	369.0	50.00
	4/15/20	600.31	3.70	596.61	0.73	ND	3.2 J	ND	200	4.4	140 JH	55	403.3	-12.50
	4/14/21	600.31	4.13	596.18	0.69	ND	ND	ND	150	1.7 J	95	70	317.4	-40.63
	7/1/21	600.31	4.80	595.71	1.6	ND	ND	0.18 J	210	3.9	150	86	453.6	-6.25
	11/19/21	600.31	3.30	597.01	0.45 J	ND	ND	ND	50	ND	73	20	143.5	-54.38
	1/12/22	600.31	3.85	596.46	1.1	ND	ND	ND	140	1.8 J	74	54	270.9	-53.75
	4/5/22	600.31	3.80	596.51	0.9	ND	ND	ND	130	1.8 J	59	75	266.7	-63.13
	7/6/22	600.31	4.11	596.20	0.73	ND	ND	ND	110	1.7 J	89	51	252.4	-44.38
	10/7/22	600.31	5.66	594.65	0.53	1.9 J	ND	ND	85	1.2 J	72	39	199.6	-55.00
	1/5/23	600.31	2.62	597.69	0.19 J	ND	ND	ND	40	ND	35	6	81.2	-78.13
	4/6/23	600.31	3.10	597.21	0.22 J	ND	ND	ND	42 J	ND	32 J	15	89.2	-80.00
	7/25/23	600.31	3.98	596.33	0.55	ND	ND	ND	89	1.3 J	90	35	215.9	-43.75
	10/3/23	600.31	6.70	593.61	0.55	ND	ND	ND	90	1.1 J	71	35	197.7	-55.63
	1/12/24	600.31	3.11	597.20	0.19 J	ND	ND	ND	36	ND	36	9.1	80.3	-77.50
	4/9/24	600.31	3.66	596.65	0.30 J	ND	ND	ND	55	0.7 J	43	22	121.0	-73.13
	7/1/24	600.31	3.27	597.04	0.18 J	ND	ND	ND	32	ND	54	11	97.2	-66.25

Notes:

1. NG = Not Gauged; NT = Not Tested; ND = Non-Detect. 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). : H = The analysis of pH was performed beyond the regulatory-required holding time of 15 minutes from the time of sample collection.
2. Water Levels measured from top of riser
3. Blue Shading = Result exceeds NY-TGS-GA for TCE
4. RED BOLD = Percent increase of TCE from Baseline
5. BLUE BOLD = Result changed as a result of data validation.
6. Data Validation was not performed on the following sample dates: 7/16/19 (sampled by others), 10/24/19 (sampled by others), 7/1/21, 11/19/21, 1/12/22.
7. 10/24/2019 data analyzed by eurolins Lancaster Laboratories Environmental, all other data analyzed by Alpha Analytical
8. QA/QC Results not included on this table, please see full analytical report.

## **ATTACHMENT C**

Well Data Sheets





## Well Data Sheet

Date: 07/01/24

Job #: 01304

Well ID: MW-3

Crew: RH + CS

Well Depth (TOR): 15.0

Well Depth (GS): 15.0

Initial Water Level (TOR): 5.26

Initial Water Level (GS): 5.86

Volume Calculation:  $(15.0 - 5.26) \times 0.163 = 1.59$

DTB-DTW\*0.163=1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity
8:54	1	7.00	1.44	20.07	0.0
9:04	2	6.99	1.40	20.06	0.0
9:14	3	7.01	1.37	20.04	0.0
9:24	4	7.02	1.35	20.02	0.0
9:35	5	7.04	1.32	20.05	0.0

Purge Method: Bailer/Submersible Pump

Initial Water Quality good

Final Water Quality good

### SAMPLE RECORD

Date: 07/01/24

Time: 9:35

Crew: RH + CS

Method: low flow

Sample ID: MW-3 (070124)

Water Quality: good

pH: 7.04

Conductivity: 1.32

Temperature: 20.05

Turbidity: 0.0

Volume: see chain

Analysis: 11

Chain of Custody #: →

Sample Type: grab

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace % 0.0

TOR= Top of Riser

GS= Ground Surface

Signature: Ryan Hooker



## Well Data Sheet

Date: 07/01/24

Job #: 01304

Well ID: MW-11

Crew: RH + CS

Well Depth (TOR): 15.05

Well Depth (GS): 15.88

Initial Water Level (TOR): 5.30

Initial Water Level (GS): 6.13

Volume Calculation:  $(15.05 - 5.30) \times 0.041 = 0.39$

DTB-DTW\*0.163=1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity
10:02	0.5	7.03	2.62	19.89	7.7
10:10	1.0	7.14	2.99	19.41	<del>68.9</del> 68.9
10:18	1.5	7.27	2.53	19.18	43.6
10:23	2.0	7.28	2.55	19.55	19.8

Purge Method: Bailer/~~Submersible Pump~~

Initial Water Quality: good

Final Water Quality: good

### SAMPLE RECORD

Date: 07/01/24

Time: 10:23

Crew: RH + CS

Method: low flow

Sample ID: MW-11 (070124)

Water Quality: good

pH: 7.28

Conductivity: 2.55

Temperature: 19.55

Turbidity: 19.8

Volume: sec chain

Analysis: "

Chain of Custody #: -

Sample Type: grab

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace @ 1.70

TOR= Top of Riser

GS= Ground Surface

Signature: Ryan Hooker



## Well Data Sheet

Date: 07/01/24

Job #: 01304

Well ID: MW-12

Crew: RH + CS

Well Depth (TOR): 14.7

Well Depth (GS): 15.2

Initial Water Level (TOR): 4.1

Initial Water Level (GS): 4.6

Volume Calculation:  $(14.7 - 4.1) \times 0.041 = 0.43$

DTB-DTW\*0.163=1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity
10:52	0.5	7.20	0.927	20.84	36.2
11:00	1.0	7.21	0.884	21.25	3.7
11:09	1.5	7.47	0.864	21.36	0.4

Purge Method: Bailer/~~Submersible Pump~~

Initial Water Quality poor

Final Water Quality good

### SAMPLE RECORD

Date: 07/01/24

Time: 11:09

Crew: RH + CS

Method: low flow

Sample ID: MW-12 (070124)

Water Quality: good

pH: 7.47

Conductivity: 0.864

Temperature: 21.36

Turbidity: 0.4

Volume: see chain

Analysis: "

Chain of Custody #: —

Sample Type: grab

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace 0.0

TOR= Top of Riser

GS= Ground Surface

Signature: Ryder Hooker



## Well Data Sheet

Date: 07/01/24

Job #: 01304

Well ID: MW-13

Crew: RH + CS

Well Depth (TOR): 14.23

Well Depth (GS): 14.93

Initial Water Level (TOR): 3.27

Initial Water Level (GS):

Volume Calculation:  $(14.23 - 3.27) \times 0.041 = 0.449$

DTB-DTW \* 0.163 = 1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity
11:37	0.5	7.25	0.594	22.58	38.9
11:43	1.0	7.30	0.551	23.21	5.3
11:50	1.5	7.36	0.546	23.52	1.9

Purge Method: Bailer/~~Submersible Pump~~

Initial Water Quality: good

Final Water Quality: good

### SAMPLE RECORD

Date: 07/01/24

Time: 11:50

Crew: RH + CS

Method: low flow

Sample ID: MW-13 (070124)

Water Quality: good

pH: 7.36

Conductivity: 0.546

Temperature: 23.52

Turbidity: 1.9

Volume: sec chain

Analysis: "

Chain of Custody #: —

Sample Type: grab

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace % 0.0

TOR= Top of Riser

GS= Ground Surface

Signature: Ryker Hooker



## Well Data Sheet

Date: 07101124

Job #: 01304

Well ID: MW - 14

Crew: RH + CS

Well Depth (TOR): 9.7

Well Depth (GS): 10.16

Initial Water Level (TOR): 6.24

Initial Water Level (GS): 6.70

Volume Calculation: No Sample

DTB-DTW\*0.163=1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity

Purge Method: Bailer/Submersible Pump

Initial Water Quality

Final Water Quality

### SAMPLE RECORD

Date:

Time:

Crew:

Method:

Sample ID:

Water Quality:

pH:

Conductivity:

Temperature:

Turbidity:

Volume:

Analysis:

Chain of Custody #:

Sample Type:

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace 0.0

TOR= Top of Riser

GS= Ground Surface

Signature: Rylee Hooker



## Well Data Sheet

Date: 07/01/24 Job #: 01304  
Well ID: MW-15  
Crew: RH + CS  
Well Depth (TOR): 10.92  
Well Depth (GS): 10.72  
Initial Water Level (TOR): 5.02  
Initial Water Level (GS): 4.82

Volume Calculation: No Sample  
DTB-DTW\*0.163=1-well vol

### Purge Record

Time	Volume	pH	Cond.	Temp.	Turbidity

Purge Method: Bailer/Submersible Pump  
Initial Water Quality  
Final Water Quality

### SAMPLE RECORD

Date: \_\_\_\_\_ Volume: \_\_\_\_\_  
Time: \_\_\_\_\_ Analysis: \_\_\_\_\_  
Crew: \_\_\_\_\_ Chain of Custody #: \_\_\_\_\_  
Method: \_\_\_\_\_ Sample Type: \_\_\_\_\_  
Sample ID: \_\_\_\_\_  
Water Quality: \_\_\_\_\_  
pH: \_\_\_\_\_  
Conductivity: \_\_\_\_\_  
Temperature: \_\_\_\_\_  
Turbidity: \_\_\_\_\_

Diameter	Multiply by
1"	0.041
2"	0.163
3"	0.367
4"	0.653
6"	1.468
8"	2.61

Comments: PID Headspace 0.0

TOR= Top of Riser  
GS= Ground Surface

Signature: Ryler Hooker



## **ATTACHMENT D**

### Analytical Laboratory Reports



## ANALYTICAL REPORT

Lab Number:	L2437281
Client:	Environmental Advantage, Inc. 3636 North Buffalo Road Orchard Park, NY 14127
ATTN:	Mark Hanna
Phone:	(716) 667-3130
Project Name:	MPC Q3 GROUNDWATER SAMPLING
Project Number:	01304
Report Date:	07/09/24

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-0826), IL (200077), IN (C-MA-03), KY (KY98045), ME (MA00086), MD (348), NJ (MA935), NY (11148), NC (25700/666), OR (MA-1316), PA (68-03671), RI (LAO00065), TX (T104704476), VT (VT-0935), VA (460195), USDA (Permit #525-23-122-91930A1).

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



**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

<b>Alpha Sample ID</b>	<b>Client ID</b>	<b>Matrix</b>	<b>Sample Location</b>	<b>Collection Date/Time</b>	<b>Receive Date</b>
L2437281-01	MW-3 (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 09:35	07/01/24
L2437281-02	MW-11 (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 10:23	07/01/24
L2437281-03	MW-11 (070124) DUPLICATE	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 10:23	07/01/24
L2437281-04	MW-12 (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 11:09	07/01/24
L2437281-05	MW-13 (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 11:50	07/01/24
L2437281-06	TRIP BLANK (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 11:38	07/01/24
L2437281-07	RINSATE BLANK (070124)	WATER	MOD-PAC CORP. BUFFALO, NY	07/01/24 12:10	07/01/24

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

### 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.

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**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**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:



Cristin Walker

Title: Technical Director/Representative

Date: 07/09/24

# ORGANICS



# **VOLATILES**

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-01      D  
**Client ID:** MW-3 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 09:35  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 23:05  
**Analyst:** MJV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westborough Lab						
Methylene chloride	ND		ug/l	6.2	1.8	2.5
1,1-Dichloroethane	ND		ug/l	6.2	1.8	2.5
Chloroform	ND		ug/l	6.2	1.8	2.5
Carbon tetrachloride	ND		ug/l	1.2	0.34	2.5
1,2-Dichloropropane	ND		ug/l	2.5	0.34	2.5
Dibromochloromethane	ND		ug/l	1.2	0.37	2.5
1,1,2-Trichloroethane	ND		ug/l	3.8	1.2	2.5
Tetrachloroethene	ND		ug/l	1.2	0.45	2.5
Chlorobenzene	ND		ug/l	6.2	1.8	2.5
Trichlorofluoromethane	ND		ug/l	6.2	1.8	2.5
1,2-Dichloroethane	ND		ug/l	1.2	0.33	2.5
1,1,1-Trichloroethane	ND		ug/l	6.2	1.8	2.5
Bromodichloromethane	ND		ug/l	1.2	0.48	2.5
trans-1,3-Dichloropropene	ND		ug/l	1.2	0.41	2.5
cis-1,3-Dichloropropene	ND		ug/l	1.2	0.36	2.5
Bromoform	ND		ug/l	5.0	1.6	2.5
1,1,2,2-Tetrachloroethane	ND		ug/l	1.2	0.42	2.5
Benzene	ND		ug/l	1.2	0.40	2.5
Toluene	ND		ug/l	6.2	1.8	2.5
Ethylbenzene	ND		ug/l	6.2	1.8	2.5
Chloromethane	ND		ug/l	6.2	1.8	2.5
Bromomethane	ND		ug/l	6.2	1.8	2.5
Vinyl chloride	1.9	J	ug/l	2.5	0.18	2.5
Chloroethane	ND		ug/l	6.2	1.8	2.5
1,1-Dichloroethene	ND		ug/l	1.2	0.42	2.5
trans-1,2-Dichloroethene	5.2	J	ug/l	6.2	1.8	2.5
Trichloroethene	250		ug/l	1.2	0.44	2.5
1,2-Dichlorobenzene	ND		ug/l	6.2	1.8	2.5

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-01 D  
**Client ID:** MW-3 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 09:35  
**Date Received:** 07/01/24  
**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	6.2	1.8	2.5
1,4-Dichlorobenzene	ND		ug/l	6.2	1.8	2.5
Methyl tert butyl ether	ND		ug/l	6.2	0.42	2.5
p/m-Xylene	ND		ug/l	6.2	1.8	2.5
o-Xylene	ND		ug/l	6.2	1.8	2.5
cis-1,2-Dichloroethene	66		ug/l	6.2	1.8	2.5
Styrene	ND		ug/l	6.2	1.8	2.5
Dichlorodifluoromethane	ND		ug/l	12	2.5	2.5
Acetone	ND		ug/l	12	3.6	2.5
Carbon disulfide	ND		ug/l	12	2.5	2.5
2-Butanone	ND		ug/l	12	4.8	2.5
4-Methyl-2-pentanone	ND		ug/l	12	2.5	2.5
2-Hexanone	ND		ug/l	12	2.5	2.5
Bromochloromethane	ND		ug/l	6.2	1.8	2.5
1,2-Dibromoethane	ND		ug/l	5.0	1.6	2.5
1,2-Dibromo-3-chloropropane	ND		ug/l	6.2	1.8	2.5
Isopropylbenzene	ND		ug/l	6.2	1.8	2.5
1,2,3-Trichlorobenzene	ND		ug/l	6.2	1.8	2.5
1,2,4-Trichlorobenzene	ND		ug/l	6.2	1.8	2.5
Methyl Acetate	ND		ug/l	5.0	0.58	2.5
Cyclohexane	ND		ug/l	25	0.68	2.5
1,4-Dioxane	ND		ug/l	620	150	2.5
Freon-113	ND		ug/l	6.2	1.8	2.5
Methyl cyclohexane	ND		ug/l	25	0.99	2.5

Surrogate	% Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	121		70-130
Toluene-d8	103		70-130
4-Bromofluorobenzene	98		70-130
Dibromofluoromethane	110		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-02  
**Client ID:** MW-11 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 10:23  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 22:15  
**Analyst:** MJV

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	0.23	J	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	14		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	0.22	J	ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	21		ug/l	2.5	0.70	1
Trichloroethene	37		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-02  
**Client ID:** MW-11 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 10:23  
**Date Received:** 07/01/24  
**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	0.17	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	17		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	3.1	J	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	117		70-130
Toluene-d8	105		70-130
4-Bromofluorobenzene	95		70-130
Dibromofluoromethane	106		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-03  
**Client ID:** MW-11 (070124) DUPLICATE  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 10:23  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 22:40  
**Analyst:** MJV

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	0.18	J	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	13		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	0.33	J	ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	20		ug/l	2.5	0.70	1
Trichloroethene	36		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-03  
**Client ID:** MW-11 (070124) DUPLICATE  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 10:23  
**Date Received:** 07/01/24  
**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	0.17	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	17		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	3.3	J	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	117		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	96		70-130
Dibromofluoromethane	112		70-130



**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-04  
**Client ID:** MW-12 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:09  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 21:50  
**Analyst:** MJV

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	0.86		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-04  
**Client ID:** MW-12 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:09  
**Date Received:** 07/01/24  
**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	0.17	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	121		70-130
Toluene-d8	101		70-130
4-Bromofluorobenzene	96		70-130
Dibromofluoromethane	111		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-05  
**Client ID:** MW-13 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:50  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 21:26  
**Analyst:** MJV

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	11		ug/l	1.0	0.07	1
Chloroethane	ND		ug/l	2.5	0.70	1
1,1-Dichloroethene	0.18	J	ug/l	0.50	0.17	1
trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1
Trichloroethene	54		ug/l	0.50	0.18	1
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-05  
**Client ID:** MW-13 (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:50  
**Date Received:** 07/01/24  
**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	0.17	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	32		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	118		70-130
Toluene-d8	99		70-130
4-Bromofluorobenzene	99		70-130
Dibromofluoromethane	109		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-06  
**Client ID:** TRIP BLANK (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:38  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 20:36  
**Analyst:** MJV

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:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-06  
**Client ID:** TRIP BLANK (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 11:38  
**Date Received:** 07/01/24  
**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	0.17	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	117		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	98		70-130
Dibromofluoromethane	109		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-07  
**Client ID:** RINSATE BLANK (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 12:10  
**Date Received:** 07/01/24  
**Field Prep:** Not Specified

**Sample Depth:**  
**Matrix:** Water  
**Analytical Method:** 1,8260D  
**Analytical Date:** 07/06/24 21:01  
**Analyst:** MJV

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:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**SAMPLE RESULTS**

**Lab ID:** L2437281-07  
**Client ID:** RINSATE BLANK (070124)  
**Sample Location:** MOD-PAC CORP. BUFFALO, NY

**Date Collected:** 07/01/24 12:10  
**Date Received:** 07/01/24  
**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	0.17	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	1.6	J	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	118		70-130
Toluene-d8	102		70-130
4-Bromofluorobenzene	98		70-130
Dibromofluoromethane	112		70-130



**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D  
 Analytical Date: 07/06/24 15:14  
 Analyst: LAC

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-07 Batch: WG1944312-5					
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:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

### Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D  
 Analytical Date: 07/06/24 15:14  
 Analyst: LAC

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-07 Batch: WG1944312-5					
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70
Methyl tert butyl ether	ND		ug/l	2.5	0.17
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:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

**Lab Number:** L2437281  
**Report Date:** 07/09/24

**Method Blank Analysis**  
**Batch Quality Control**

Analytical Method: 1,8260D  
 Analytical Date: 07/06/24 15:14  
 Analyst: LAC

Parameter	Result	Qualifier	Units	RL	MDL
Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-07 Batch: WG1944312-5					

Surrogate	%Recovery	Qualifier	Acceptance Criteria
1,2-Dichloroethane-d4	114		70-130
Toluene-d8	101		70-130
4-Bromofluorobenzene	100		70-130
Dibromofluoromethane	109		70-130

# **Lab Control Sample Analysis** Batch Quality Control

**Project Name:** MPC Q3 GROUNDWATER SAMPLING

**Project Number:** 01304

**Lab Number:** L2437281

**Report Date:** 07/09/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 Batch: WG1944312-3 WG1944312-4								
Methylene chloride	92		92		70-130	0		20
1,1-Dichloroethane	100		100		70-130	0		20
Chloroform	97		95		70-130	2		20
Carbon tetrachloride	90		90		63-132	0		20
1,2-Dichloropropane	98		100		70-130	2		20
Dibromochloromethane	93		97		63-130	4		20
1,1,2-Trichloroethane	100		100		70-130	0		20
Tetrachloroethene	91		53	Q	70-130	53	Q	20
Chlorobenzene	99		98		75-130	1		20
Trichlorofluoromethane	92		92		62-150	0		20
1,2-Dichloroethane	98		99		70-130	1		20
1,1,1-Trichloroethane	91		91		67-130	0		20
Bromodichloromethane	96		96		67-130	0		20
trans-1,3-Dichloropropene	96		96		70-130	0		20
cis-1,3-Dichloropropene	96		96		70-130	0		20
Bromoform	94		95		54-136	1		20
1,1,2,2-Tetrachloroethane	110		120		67-130	9		20
Benzene	98		98		70-130	0		20
Toluene	96		98		70-130	2		20
Ethylbenzene	97		97		70-130	0		20
Chloromethane	76		76		64-130	0		20
Bromomethane	45		44		39-139	2		20
Vinyl chloride	90		88		55-140	2		20

# **Lab Control Sample Analysis** **Batch Quality Control**

**Project Name:** MPC Q3 GROUNDWATER SAMPLING

**Project Number:** 01304

**Lab Number:** L2437281

**Report Date:** 07/09/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 Batch: WG1944312-3 WG1944312-4								
Chloroethane	110		100		55-138	10		20
1,1-Dichloroethene	83		86		61-145	4		20
trans-1,2-Dichloroethene	90		92		70-130	2		20
Trichloroethene	82		83		70-130	1		20
1,2-Dichlorobenzene	96		96		70-130	0		20
1,3-Dichlorobenzene	100		98		70-130	2		20
1,4-Dichlorobenzene	99		98		70-130	1		20
Methyl tert butyl ether	92		95		63-130	3		20
p/m-Xylene	95		100		70-130	5		20
o-Xylene	100		100		70-130	0		20
cis-1,2-Dichloroethene	94		96		70-130	2		20
Styrene	100		100		70-130	0		20
Dichlorodifluoromethane	76		75		36-147	1		20
Acetone	87		92		58-148	6		20
Carbon disulfide	96		89		51-130	8		20
2-Butanone	85		93		63-138	9		20
4-Methyl-2-pentanone	91		97		59-130	6		20
2-Hexanone	84		92		57-130	9		20
Bromochloromethane	100		100		70-130	0		20
1,2-Dibromoethane	92		94		70-130	2		20
1,2-Dibromo-3-chloropropane	86		88		41-144	2		20
Isopropylbenzene	95		93		70-130	2		20
1,2,3-Trichlorobenzene	88		90		70-130	2		20

# Lab Control Sample Analysis

## Batch Quality Control

**Project Name:** MPC Q3 GROUNDWATER SAMPLING

**Project Number:** 01304

**Lab Number:** L2437281

**Report Date:** 07/09/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 Batch: WG1944312-3 WG1944312-4								
1,2,4-Trichlorobenzene	90		90		70-130	0		20
Methyl Acetate	100		100		70-130	0		20
Cyclohexane	90		91		70-130	1		20
1,4-Dioxane	84		90		56-162	7		20
Freon-113	87		86		70-130	1		20
Methyl cyclohexane	86		86		70-130	0		20

Surrogate	LCS %Recovery	Qual	LCSD %Recovery	Qual	Acceptance Criteria
1,2-Dichloroethane-d4	105		110		70-130
Toluene-d8	102		103		70-130
4-Bromofluorobenzene	99		97		70-130
Dibromofluoromethane	101		102		70-130

# Matrix Spike Analysis

## Batch Quality Control

**Project Name:** MPC Q3 GROUNDWATER SAMPLING

**Project Number:** 01304

**Lab Number:** L2437281

**Report Date:** 07/09/24

Parameter	Native Sample	MS Added	MS Found	MS %Recovery	Qual	MSD Found	MSD %Recovery	Qual	Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 QC Batch ID: WG1944312-8 WG1944312-9 QC Sample: L2437281-04 Client ID: MW-12 (070124)												
Methylene chloride	ND	10	11	110		10	100		70-130	10		20
1,1-Dichloroethane	ND	10	12	120		12	120		70-130	0		20
Chloroform	ND	10	11	110		11	110		70-130	0		20
Carbon tetrachloride	ND	10	12	120		11	110		63-132	9		20
1,2-Dichloropropane	ND	10	12	120		11	110		70-130	9		20
Dibromochloromethane	ND	10	11	110		10	100		63-130	10		20
1,1,2-Trichloroethane	ND	10	11	110		11	110		70-130	0		20
Tetrachloroethene	ND	10	10	100		10	100		70-130	0		20
Chlorobenzene	ND	10	11	110		10	100		75-130	10		20
Trichlorofluoromethane	ND	10	13	130		13	130		62-150	0		20
1,2-Dichloroethane	ND	10	12	120		12	120		70-130	0		20
1,1,1-Trichloroethane	ND	10	12	120		11	110		67-130	9		20
Bromodichloromethane	ND	10	11	110		11	110		67-130	0		20
trans-1,3-Dichloropropene	ND	10	10	100		9.8	98		70-130	2		20
cis-1,3-Dichloropropene	ND	10	10	100		9.7	97		70-130	3		20
Bromoform	ND	10	10	100		9.7	97		54-136	3		20
1,1,2,2-Tetrachloroethane	ND	10	12	120		12	120		67-130	0		20
Benzene	ND	10	12	120		11	110		70-130	9		20
Toluene	ND	10	11	110		11	110		70-130	0		20
Ethylbenzene	ND	10	11	110		11	110		70-130	0		20
Chloromethane	ND	10	11	110		11	110		64-130	0		20
Bromomethane	ND	10	5.8	58		5.8	58		39-139	0		20
Vinyl chloride	ND	10	12	120		12	120		55-140	0		20

**Matrix Spike Analysis****Batch Quality Control****Project Name:** MPC Q3 GROUNDWATER SAMPLING**Project Number:** 01304**Lab Number:** L2437281**Report Date:** 07/09/24

<b>Parameter</b>	<b>Native Sample</b>	<b>MS Added</b>	<b>MS Found</b>	<b>MS %Recovery</b>	<b>Qual</b>	<b>MSD Found</b>	<b>MSD %Recovery</b>	<b>Qual</b>	<b>Recovery Limits</b>	<b>RPD</b>	<b>Qual</b>	<b>RPD Limits</b>
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 QC Batch ID: WG1944312-8 WG1944312-9 QC Sample: L2437281-04 Client ID: MW-12 (070124)												
Chloroethane	ND	10	13	130		14	140	Q	55-138	7		20
1,1-Dichloroethene	ND	10	11	110		10	100		61-145	10		20
trans-1,2-Dichloroethene	ND	10	11	110		10	100		70-130	10		20
Trichloroethene	0.86	10	11	101		10	91		70-130	10		20
1,2-Dichlorobenzene	ND	10	10	100		10	100		70-130	0		20
1,3-Dichlorobenzene	ND	10	10	100		10	100		70-130	0		20
1,4-Dichlorobenzene	ND	10	10	100		10	100		70-130	0		20
Methyl tert butyl ether	ND	10	10	100		9.7	97		63-130	3		20
p/m-Xylene	ND	20	23	115		22	110		70-130	4		20
o-Xylene	ND	20	23	115		22	110		70-130	4		20
cis-1,2-Dichloroethene	ND	10	11	110		10	100		70-130	10		20
Styrene	ND	20	23	115		22	110		70-130	4		20
Dichlorodifluoromethane	ND	10	10	100		10	100		36-147	0		20
Acetone	ND	10	12	120		13	130		58-148	8		20
Carbon disulfide	ND	10	11	110		11	110		51-130	0		20
2-Butanone	ND	10	12	120		12	120		63-138	0		20
4-Methyl-2-pentanone	ND	10	11	110		10	100		59-130	10		20
2-Hexanone	ND	10	10	100		9.8	98		57-130	2		20
Bromochloromethane	ND	10	11	110		11	110		70-130	0		20
1,2-Dibromoethane	ND	10	11	110		10	100		70-130	10		20
1,2-Dibromo-3-chloropropane	ND	10	9.4	94		8.9	89		41-144	5		20
Isopropylbenzene	ND	10	10	100		10	100		70-130	0		20
1,2,3-Trichlorobenzene	ND	10	9.3	93		9.3	93		70-130	0		20



**Matrix Spike Analysis***Batch Quality Control***Project Name:** MPC Q3 GROUNDWATER SAMPLING**Project Number:** 01304**Lab Number:** L2437281**Report Date:** 07/09/24

<b>Parameter</b>	<b>Native Sample</b>	<b>MS Added</b>	<b>MS Found</b>	<b>MS %Recovery</b>	<b>Qual</b>	<b>MSD Found</b>	<b>MSD %Recovery</b>	<b>Qual</b>	<b>Recovery Limits</b>	<b>RPD</b>	<b>Qual</b>	<b>RPD Limits</b>
Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-07 QC Batch ID: WG1944312-8 WG1944312-9 QC Sample: L2437281-04 Client ID: MW-12 (070124)												
1,2,4-Trichlorobenzene	ND	10	9.0	90		9.1	91		70-130	1		20
Methyl Acetate	ND	10	11	110		11	110		70-130	0		20
Cyclohexane	ND	10	13	130		12	120		70-130	8		20
1,4-Dioxane	ND	500	460	92		470	94		56-162	2		20
Freon-113	ND	10	11	110		10	100		70-130	10		20
Methyl cyclohexane	ND	10	10	100		10	100		70-130	0		20

<b>Surrogate</b>	<b>MS % Recovery</b>	<b>MS Qualifier</b>	<b>MSD % Recovery</b>	<b>MSD Qualifier</b>	<b>Acceptance Criteria</b>
1,2-Dichloroethane-d4	115		114		70-130
4-Bromofluorobenzene	97		97		70-130
Dibromofluoromethane	112		104		70-130
Toluene-d8	103		102		70-130

**Project Name:** MPC Q3 GROUNDWATER SAMPLING**Lab Number:** L2437281**Project Number:** 01304**Report Date:** 07/09/24**Sample Receipt and Container Information**

Were project specific reporting limits specified?

YES

**Cooler Information**

<b>Cooler</b>	<b>Custody Seal</b>
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A	Absent
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B	Absent
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C	Absent
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**Container Information**

<b>Container ID</b>	<b>Container Type</b>	<b>Cooler</b>	<b>Initial pH</b>	<b>Final pH</b>	<b>Temp deg C</b>	<b>Pres</b>	<b>Seal</b>	<b>Frozen Date/Time</b>	<b>Analysis(*)</b>
L2437281-01A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-01B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-01C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-02A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-02B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-02C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-03A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-03B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-03C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04A1	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04A2	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04B1	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04B2	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04C1	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-04C2	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-05A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-05B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-05C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
**Project Number:** 01304

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**Container Information**

<b>Container ID</b>	<b>Container Type</b>	<b>Cooler</b>	<b>Initial pH</b>	<b>Final pH</b>	<b>Temp deg C</b>	<b>Pres</b>	<b>Seal</b>	<b>Frozen Date/Time</b>	<b>Analysis(*)</b>
L2437281-06A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-06B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-07A	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-07B	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)
L2437281-07C	Vial HCl preserved	A	NA		2.9	Y	Absent		NYTCL-8260-R2(14)

**Project Name:** MPC Q3 GROUNDWATER SAMPLING  
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**Report Date:** 07/09/24

## 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.

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### 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.

**Chlordane:** 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.)

**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'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**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 for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**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

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#### Data Qualifiers

Identified Compounds (TICs). For calculated parameters, this represents that one or more values used in the calculation were estimated.

- 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 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.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

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## 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.



**Alpha Analytical, Inc.**Facility: **Company-wide**Department: **Quality Assurance**Title: **Certificate/Approval Program Summary**ID No.: **17873**

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**Certification Information**

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

**Westborough Facility****EPA 624.1:** m/p-xylene, o-xylene, Naphthalene**EPA 625.1:** alpha-Terpineol**EPA 8260D:** NPW: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; SCM: Iodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene.**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.**Mansfield Facility****SM 2540D:** TSS.**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.

**Nonpotable Water:** **EPA RSK-175 Dissolved 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 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).**Microbiology:** **SM9223B-Colilert-QT; Enterolert-QT, 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.



 <b>NEW YORK CHAIN OF CUSTODY</b> Westborough, MA 01581 8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193 Mansfield, MA 02048 320 Forbes Blvd TEL: 508-822-9300 FAX: 508-822-3286		<b>Service Centers</b> Mahwah, NJ 07430: 35 Whitney Rd, Suite 5 Albany, NY 12205: 14 Walker Way Tonawanda, NY 14150: 275 Cooper Ave, Suite 105		Page 1 of 1		Date Rec'd in Lab <b>7/02/24</b>		ALPHA Job # <b>L2437281</b>																																																																																																																																																																																																							
		<b>Project Information</b> Project Name: <b>MPC g3 Groundwater Sampling</b> Project Location: <b>Mod-Pac Corp. Buffalo, NY</b> Project # <b>01304</b> (Use Project name as Project #) <input type="checkbox"/> Project Manager: <b>Mark Hanna + Mary Szustak</b> ALPHAQuote #: Turn-Around Time Standard <input checked="" type="checkbox"/> Rush (only if pre approved) <input type="checkbox"/> Due Date: # of Days:		<b>Deliverables</b> <input type="checkbox"/> ASP-A <input checked="" type="checkbox"/> ASP-B <input type="checkbox"/> EQUIS (1 File) <input checked="" type="checkbox"/> EQUIS (4 File) <input type="checkbox"/> Other		<b>Billing Information</b> <input checked="" type="checkbox"/> Same as Client Info PO # <b>01304</b>																																																																																																																																																																																																									
<b>Client Information</b> Client: <b>Env. Advantage, Inc.</b> Address: <b>3636 N. Buffalo Rd</b> <b>Orchard Park, NY 14127</b> Phone: <b>(716) 667-3130</b> Fax: <b>(716) 667-3156</b> Email: <b>mhanna@envadvantage.com</b>		<b>Regulatory Requirement</b> <input type="checkbox"/> NY TOGS <input type="checkbox"/> NY Part 375 <input type="checkbox"/> AWQ Standards <input type="checkbox"/> NY CP-51 <input type="checkbox"/> NY Restricted Use <input type="checkbox"/> Other <input type="checkbox"/> NY Unrestricted Use <input type="checkbox"/> NYC Sewer Discharge		<b>Disposal Site Information</b> Please identify below location of applicable disposal facilities. Disposal Facility: <input type="checkbox"/> NJ <input type="checkbox"/> NY <input type="checkbox"/> Other:																																																																																																																																																																																																											
These samples have been previously analyzed by Alpha <input type="checkbox"/> Other project specific requirements/comments: <b>Please also email results to labresults@envadvantage.com</b> Please specify Metals or TAL.		<b>ANALYSIS</b> VOCs 8260 TCL		<b>Sample Filtration</b> <input type="checkbox"/> Done <input type="checkbox"/> Lab to do <input type="checkbox"/> Preservation <input type="checkbox"/> Lab to do (Please Specify below)		<b>Sample Specific Comments</b>																																																																																																																																																																																																									
<table border="1" style="width:100%; border-collapse: collapse;"> <thead> <tr> <th rowspan="2">ALPHA Lab ID (Lab Use Only)</th> <th rowspan="2">Sample ID</th> <th colspan="2">Collection</th> <th rowspan="2">Sample Matrix</th> <th rowspan="2">Sampler's Initials</th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> <th rowspan="2"></th> </tr> <tr> <th>Date</th> <th>Time</th> </tr> </thead> <tbody> <tr> <td>37281-01</td> <td>MW-3 (070124)</td> <td>07/01/24</td> <td>9:35</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-02</td> <td>MW-11 (070124)</td> <td>07/01/24</td> <td>10:23</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-03</td> <td>MW-11 (070124) duplicate</td> <td>07/01/24</td> <td>10:23</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-04</td> <td>MW-12 (070124)</td> <td>07/01/24</td> <td>11:09</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-04</td> <td>MW-12 (070124) MS</td> <td>07/01/24</td> <td>11:09</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-04</td> <td>MW-12 (070124) MSD</td> <td>07/01/24</td> <td>11:09</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>05</td> <td>MW-13 (070124)</td> <td>07/01/24</td> <td>11:50</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>06</td> <td>Trip Blank (070124)</td> <td>07/01/24</td> <td>11:38</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> <tr> <td>-07</td> <td>Rinsate Blank (070124)</td> <td>07/01/24</td> <td>12:10</td> <td>Wa</td> <td>RH</td> <td>X</td> <td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td><td></td> </tr> </tbody> </table>		ALPHA Lab ID (Lab Use Only)	Sample ID	Collection		Sample Matrix	Sampler's Initials															Date	Time	37281-01	MW-3 (070124)	07/01/24	9:35	Wa	RH	X														-02	MW-11 (070124)	07/01/24	10:23	Wa	RH	X														-03	MW-11 (070124) duplicate	07/01/24	10:23	Wa	RH	X														-04	MW-12 (070124)	07/01/24	11:09	Wa	RH	X														-04	MW-12 (070124) MS	07/01/24	11:09	Wa	RH	X														-04	MW-12 (070124) MSD	07/01/24	11:09	Wa	RH	X														05	MW-13 (070124)	07/01/24	11:50	Wa	RH	X														06	Trip Blank (070124)	07/01/24	11:38	Wa	RH	X														-07	Rinsate Blank (070124)	07/01/24	12:10	Wa	RH	X														Preservative Code: A = None B = HCl C = HNO <sub>3</sub> D = H <sub>2</sub> SO <sub>4</sub> E = NaOH F = MeOH G = NaHSO <sub>4</sub> H = Na <sub>2</sub> S <sub>2</sub> O <sub>3</sub> K/E = Zn Ac/NaOH O = Other Container Code: P = Plastic A = Amber Glass V = Vial G = Glass B = Bacteria Cup C = Cube O = Other E = Encore D = BOD Bottle Westboro: Certification No: MA935 Mansfield: Certification No: MA015 Container Type <b>V</b> Preservative <b>B</b>		Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are resolved. BY EXECUTING THIS COC, THE CLIENT HAS READ AND AGREES TO BE BOUND BY ALPHA'S TERMS & CONDITIONS. (See reverse side.)	
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Form No: 01-25 HC (rev. 30-Sept-2013)		Relinquished By: <b>T. Balkin</b> Date/Time: <b>7/1/24 15:40</b>		Received By: <b>T. Balkin</b> Date/Time: <b>7-1-24 15:40</b>		Relinquished By: <b>M. Hanna</b> Date/Time: <b>7-1-24 16:30</b>		Received By: <b>M. Hanna</b> Date/Time: <b>7-1-24 16:30</b>																																																																																																																																																																																																							



## ANALYTICAL REPORT

Lab Number:	L2453073
Client:	Environmental Advantage, Inc. 3636 North Buffalo Road Orchard Park, NY 14127
ATTN:	Mark Hanna
Phone:	(716) 667-3130
Project Name:	Q3 2024 SSDS MONITORING
Project Number:	01304
Report Date:	09/24/24

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Certifications & Approvals: MA (M-MA030), NH NELAP (2062), CT (PH-0825), DoD (L2474), FL (E87814), IL (200081), IN (C-MA-04), KY (KY98046), LA (85084), ME (MA00030), MD (350), MI (9110), MN (025-999-495), NJ (MA015), NY (11627), NC (685), OR (MA-0262), PA (68-02089), RI (LAO00299), TX (T104704419), VT (VT-0015), VA (460194), WA (C954), US Army Corps of Engineers, USDA (Permit #525-23-107-88708A1), USFWS (Permit #A24920).

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320 Forbes Boulevard, Mansfield, MA 02048-1806  
508-822-9300 (Fax) 508-822-3288 800-624-9220 - [www.alphalab.com](http://www.alphalab.com)



**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

<b>Alpha Sample ID</b>	<b>Client ID</b>	<b>Matrix</b>	<b>Sample Location</b>	<b>Collection Date/Time</b>	<b>Receive Date</b>
L2453073-01	AREA A-POST(091624)	SOIL_VAPOR	1801 ELMWOOD AVE.	09/16/24 11:50	09/16/24
L2453073-02	AREA A-PRE(091624)	SOIL_VAPOR	1801 ELMWOOD AVE.	09/16/24 11:55	09/16/24

**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

### 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 and solids are reported on a dry weight basis unless otherwise noted. Tissues are reported "as received" or on a wet 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:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

**Case Narrative (continued)**

Volatile Organics in Air

L2453073-01 and -02: Samples were transferred from a Tedlar bag into a fused silica lined canister upon receipt in order to extend the holding time for analysis.

L2453073-02D: The sample has elevated detection limits due to the dilution required by the elevated concentrations of target compounds in the sample.

The WG1975275-3 LCS recovery for 1,3-butadiene (140%), n-hexane (140%), 1,1,1-trichloroethane (135%), cyclohexane (142%), bromodichloromethane (150%), 1,4-dioxane (146%), 2,2,4-trimethylpentane (138%), and sec-butylbenzene (69%), associated with , is above the upper 130% acceptance limit. All samples associated with this LCS do not have reportable amounts of this analyte.

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:  Christopher J. Anderson

Title: Technical Director/Representative

Date: 09/24/24

**AIR**

**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-01  
 Client ID: AREA A-POST(091624)  
 Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:50  
 Date Received: 09/16/24  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Soil\_Vapor  
 Analytical Method: 48,TO-15  
 Analytical Date: 09/22/24 23:09  
 Analyst: KJD

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
Dichlorodifluoromethane	0.470	0.200	--	2.32	0.989	--		1
Chloromethane	0.469	0.200	--	0.968	0.413	--		1
Freon-114	ND	0.200	--	ND	1.40	--		1
Vinyl chloride	ND	0.200	--	ND	0.511	--		1
1,3-Butadiene	ND	0.200	--	ND	0.442	--		1
Bromomethane	ND	0.200	--	ND	0.777	--		1
Chloroethane	ND	0.200	--	ND	0.528	--		1
Ethanol	33.3	5.00	--	62.7	9.42	--		1
Vinyl bromide	ND	0.200	--	ND	0.874	--		1
Acetone	13.7	1.00	--	32.5	2.38	--		1
Trichlorofluoromethane	0.648	0.200	--	3.64	1.12	--		1
Isopropanol	93.2	0.500	--	229	1.23	--		1
1,1-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Tertiary butyl Alcohol	ND	0.500	--	ND	1.52	--		1
Methylene chloride	1.05	0.500	--	3.65	1.74	--		1
3-Chloropropene	ND	0.200	--	ND	0.626	--		1
Carbon disulfide	3.30	0.200	--	10.3	0.623	--		1
Freon-113	ND	0.200	--	ND	1.53	--		1
trans-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
1,1-Dichloroethane	ND	0.200	--	ND	0.809	--		1
Methyl tert butyl ether	ND	0.200	--	ND	0.721	--		1
2-Butanone	ND	0.500	--	ND	1.47	--		1
cis-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1



**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-01  
 Client ID: AREA A-POST(091624)  
 Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:50  
 Date Received: 09/16/24  
 Field Prep: Not Specified

Sample Depth:

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
Ethyl Acetate	ND	0.500	--	ND	1.80	--		1
Chloroform	ND	0.200	--	ND	0.977	--		1
Tetrahydrofuran	ND	0.500	--	ND	1.47	--		1
1,2-Dichloroethane	ND	0.200	--	ND	0.809	--		1
n-Hexane	3.74	0.200	--	13.2	0.705	--		1
1,1,1-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Benzene	1.10	0.200	--	3.51	0.639	--		1
Carbon tetrachloride	ND	0.200	--	ND	1.26	--		1
Cyclohexane	0.300	0.200	--	1.03	0.688	--		1
1,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
Bromodichloromethane	ND	0.200	--	ND	1.34	--		1
1,4-Dioxane	ND	0.200	--	ND	0.721	--		1
Trichloroethene	ND	0.200	--	ND	1.07	--		1
2,2,4-Trimethylpentane	0.345	0.200	--	1.61	0.934	--		1
Heptane	0.672	0.200	--	2.75	0.820	--		1
cis-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
4-Methyl-2-pentanone	1.58	0.500	--	6.48	2.05	--		1
trans-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
1,1,2-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Toluene	6.28	0.200	--	23.7	0.754	--		1
2-Hexanone	ND	0.200	--	ND	0.820	--		1
Dibromochloromethane	ND	0.200	--	ND	1.70	--		1
1,2-Dibromoethane	ND	0.200	--	ND	1.54	--		1
Tetrachloroethene	ND	0.200	--	ND	1.36	--		1
Chlorobenzene	ND	0.200	--	ND	0.921	--		1
Ethylbenzene	2.42	0.200	--	10.5	0.869	--		1





**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-01

Client ID: AREA A-POST(091624)

Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:50

Date Received: 09/16/24

Field Prep: Not Specified

Sample Depth:

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
p/m-Xylene	11.8	0.400	--	51.3	1.74	--		1
Bromoform	ND	0.200	--	ND	2.07	--		1
Styrene	ND	0.200	--	ND	0.852	--		1
1,1,2,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
o-Xylene	3.77	0.200	--	16.4	0.869	--		1
4-Ethyltoluene	0.595	0.200	--	2.93	0.983	--		1
1,3,5-Trimethylbenzene	0.733	0.200	--	3.60	0.983	--		1
1,2,4-Trimethylbenzene	2.92	0.200	--	14.4	0.983	--		1
Benzyl chloride	ND	0.200	--	ND	1.04	--		1
1,3-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,4-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2,4-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Naphthalene	ND	0.200	--	ND	1.05	--		1
Hexachlorobutadiene	ND	0.200	--	ND	2.13	--		1

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	97		60-140
Bromochloromethane	98		60-140
chlorobenzene-d5	98		60-140



**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-02 D  
 Client ID: AREA A-PRE(091624)  
 Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:55  
 Date Received: 09/16/24  
 Field Prep: Not Specified

Sample Depth:  
 Matrix: Soil\_Vapor  
 Analytical Method: 48,TO-15  
 Analytical Date: 09/23/24 07:15  
 Analyst: KJD

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
Dichlorodifluoromethane	0.478	0.476	--	2.36	2.35	--		2.381
Chloromethane	ND	0.476	--	ND	0.983	--		2.381
Freon-114	ND	0.476	--	ND	3.33	--		2.381
Vinyl chloride	ND	0.476	--	ND	1.22	--		2.381
1,3-Butadiene	ND	0.476	--	ND	1.05	--		2.381
Bromomethane	ND	0.476	--	ND	1.85	--		2.381
Chloroethane	ND	0.476	--	ND	1.26	--		2.381
Ethanol	31.7	11.9	--	59.7	22.4	--		2.381
Vinyl bromide	ND	0.476	--	ND	2.08	--		2.381
Acetone	82.1	2.38	--	195	5.65	--		2.381
Trichlorofluoromethane	0.998	0.476	--	5.61	2.67	--		2.381
Isopropanol	455	1.19	--	1120	2.93	--		2.381
1,1-Dichloroethene	ND	0.476	--	ND	1.89	--		2.381
Tertiary butyl Alcohol	3.61	1.19	--	10.9	3.61	--		2.381
Methylene chloride	ND	1.19	--	ND	4.13	--		2.381
3-Chloropropene	ND	0.476	--	ND	1.49	--		2.381
Carbon disulfide	1.18	0.476	--	3.67	1.48	--		2.381
Freon-113	ND	0.476	--	ND	3.65	--		2.381
trans-1,2-Dichloroethene	ND	0.476	--	ND	1.89	--		2.381
1,1-Dichloroethane	ND	0.476	--	ND	1.93	--		2.381
Methyl tert butyl ether	ND	0.476	--	ND	1.72	--		2.381
2-Butanone	1.48	1.19	--	4.36	3.51	--		2.381
cis-1,2-Dichloroethene	1.08	0.476	--	4.28	1.89	--		2.381



**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-02 D  
 Client ID: AREA A-PRE(091624)  
 Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:55  
 Date Received: 09/16/24  
 Field Prep: Not Specified

Sample Depth:

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
Ethyl Acetate	ND	1.19	--	ND	4.29	--		2.381
Chloroform	0.514	0.476	--	2.51	2.32	--		2.381
Tetrahydrofuran	ND	1.19	--	ND	3.51	--		2.381
1,2-Dichloroethane	ND	0.476	--	ND	1.93	--		2.381
n-Hexane	7.37	0.476	--	26.0	1.68	--		2.381
1,1,1-Trichloroethane	ND	0.476	--	ND	2.60	--		2.381
Benzene	0.683	0.476	--	2.18	1.52	--		2.381
Carbon tetrachloride	ND	0.476	--	ND	2.99	--		2.381
Cyclohexane	1.78	0.476	--	6.13	1.64	--		2.381
1,2-Dichloropropane	ND	0.476	--	ND	2.20	--		2.381
Bromodichloromethane	ND	0.476	--	ND	3.19	--		2.381
1,4-Dioxane	ND	0.476	--	ND	1.72	--		2.381
Trichloroethene	74.4	0.476	--	400	2.56	--		2.381
2,2,4-Trimethylpentane	0.540	0.476	--	2.52	2.22	--		2.381
Heptane	2.79	0.476	--	11.4	1.95	--		2.381
cis-1,3-Dichloropropene	ND	0.476	--	ND	2.16	--		2.381
4-Methyl-2-pentanone	2.01	1.19	--	8.24	4.88	--		2.381
trans-1,3-Dichloropropene	ND	0.476	--	ND	2.16	--		2.381
1,1,2-Trichloroethane	ND	0.476	--	ND	2.60	--		2.381
Toluene	5.84	0.476	--	22.0	1.79	--		2.381
2-Hexanone	ND	0.476	--	ND	1.95	--		2.381
Dibromochloromethane	ND	0.476	--	ND	4.06	--		2.381
1,2-Dibromoethane	ND	0.476	--	ND	3.66	--		2.381
Tetrachloroethene	ND	0.476	--	ND	3.23	--		2.381
Chlorobenzene	ND	0.476	--	ND	2.19	--		2.381
Ethylbenzene	2.23	0.476	--	9.69	2.07	--		2.381



**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24**SAMPLE RESULTS**

Lab ID: L2453073-02 D  
 Client ID: AREA A-PRE(091624)  
 Sample Location: 1801 ELMWOOD AVE.

Date Collected: 09/16/24 11:55  
 Date Received: 09/16/24  
 Field Prep: Not Specified

Sample Depth:

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab								
p/m-Xylene	10.1	0.952	--	43.9	4.14	--		2.381
Bromoform	ND	0.476	--	ND	4.92	--		2.381
Styrene	ND	0.476	--	ND	2.03	--		2.381
1,1,2,2-Tetrachloroethane	ND	0.476	--	ND	3.27	--		2.381
o-Xylene	3.32	0.476	--	14.4	2.07	--		2.381
4-Ethyltoluene	1.77	0.476	--	8.70	2.34	--		2.381
1,3,5-Trimethylbenzene	0.860	0.476	--	4.23	2.34	--		2.381
1,2,4-Trimethylbenzene	3.33	0.476	--	16.4	2.34	--		2.381
Benzyl chloride	ND	0.476	--	ND	2.46	--		2.381
1,3-Dichlorobenzene	ND	0.476	--	ND	2.86	--		2.381
1,4-Dichlorobenzene	ND	0.476	--	ND	2.86	--		2.381
1,2-Dichlorobenzene	ND	0.476	--	ND	2.86	--		2.381
1,2,4-Trichlorobenzene	ND	0.476	--	ND	3.53	--		2.381
Naphthalene	ND	0.476	--	ND	2.50	--		2.381
Hexachlorobutadiene	ND	0.476	--	ND	5.08	--		2.381

Internal Standard	% Recovery	Qualifier	Acceptance Criteria
1,4-Difluorobenzene	95		60-140
Bromochloromethane	95		60-140
chlorobenzene-d5	99		60-140



Project Name: Q3 2024 SSDS MONITORING

Lab Number: L2453073

Project Number: 01304

Report Date: 09/24/24

### Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 09/22/24 16:40

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab for sample(s): 01-02 Batch: WG1974841-4								
Dichlorodifluoromethane	ND	0.200	--	ND	0.989	--		1
Chloromethane	ND	0.200	--	ND	0.413	--		1
Freon-114	ND	0.200	--	ND	1.40	--		1
Vinyl chloride	ND	0.200	--	ND	0.511	--		1
1,3-Butadiene	ND	0.200	--	ND	0.442	--		1
Bromomethane	ND	0.200	--	ND	0.777	--		1
Chloroethane	ND	0.200	--	ND	0.528	--		1
Ethanol	ND	5.00	--	ND	9.42	--		1
Vinyl bromide	ND	0.200	--	ND	0.874	--		1
Acetone	ND	1.00	--	ND	2.38	--		1
Trichlorofluoromethane	ND	0.200	--	ND	1.12	--		1
Isopropanol	ND	0.500	--	ND	1.23	--		1
1,1-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Tertiary butyl Alcohol	ND	0.500	--	ND	1.52	--		1
Methylene chloride	ND	0.500	--	ND	1.74	--		1
3-Chloropropene	ND	0.200	--	ND	0.626	--		1
Carbon disulfide	ND	0.200	--	ND	0.623	--		1
Freon-113	ND	0.200	--	ND	1.53	--		1
trans-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
1,1-Dichloroethane	ND	0.200	--	ND	0.809	--		1
Methyl tert butyl ether	ND	0.200	--	ND	0.721	--		1
2-Butanone	ND	0.500	--	ND	1.47	--		1
cis-1,2-Dichloroethene	ND	0.200	--	ND	0.793	--		1
Ethyl Acetate	ND	0.500	--	ND	1.80	--		1
Chloroform	ND	0.200	--	ND	0.977	--		1



Project Name: Q3 2024 SSDS MONITORING

Lab Number: L2453073

Project Number: 01304

Report Date: 09/24/24

### Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 09/22/24 16:40

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab for sample(s): 01-02 Batch: WG1974841-4								
Tetrahydrofuran	ND	0.500	--	ND	1.47	--		1
1,2-Dichloroethane	ND	0.200	--	ND	0.809	--		1
n-Hexane	ND	0.200	--	ND	0.705	--		1
1,1,1-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Benzene	ND	0.200	--	ND	0.639	--		1
Carbon tetrachloride	ND	0.200	--	ND	1.26	--		1
Cyclohexane	ND	0.200	--	ND	0.688	--		1
1,2-Dichloropropane	ND	0.200	--	ND	0.924	--		1
Bromodichloromethane	ND	0.200	--	ND	1.34	--		1
1,4-Dioxane	ND	0.200	--	ND	0.721	--		1
Trichloroethene	ND	0.200	--	ND	1.07	--		1
2,2,4-Trimethylpentane	ND	0.200	--	ND	0.934	--		1
Heptane	ND	0.200	--	ND	0.820	--		1
cis-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
4-Methyl-2-pentanone	ND	0.500	--	ND	2.05	--		1
trans-1,3-Dichloropropene	ND	0.200	--	ND	0.908	--		1
1,1,2-Trichloroethane	ND	0.200	--	ND	1.09	--		1
Toluene	ND	0.200	--	ND	0.754	--		1
2-Hexanone	ND	0.200	--	ND	0.820	--		1
Dibromochloromethane	ND	0.200	--	ND	1.70	--		1
1,2-Dibromoethane	ND	0.200	--	ND	1.54	--		1
Tetrachloroethene	ND	0.200	--	ND	1.36	--		1
Chlorobenzene	ND	0.200	--	ND	0.921	--		1
Ethylbenzene	ND	0.200	--	ND	0.869	--		1
p/m-Xylene	ND	0.400	--	ND	1.74	--		1



Project Name: Q3 2024 SSDS MONITORING

Lab Number: L2453073

Project Number: 01304

Report Date: 09/24/24

### Method Blank Analysis Batch Quality Control

Analytical Method: 48,TO-15

Analytical Date: 09/22/24 16:40

Parameter	ppbV			ug/m3			Qualifier	Dilution Factor
	Results	RL	MDL	Results	RL	MDL		
Volatile Organics in Air - Mansfield Lab for sample(s): 01-02 Batch: WG1974841-4								
Bromoform	ND	0.200	--	ND	2.07	--		1
Styrene	ND	0.200	--	ND	0.852	--		1
1,1,2,2-Tetrachloroethane	ND	0.200	--	ND	1.37	--		1
o-Xylene	ND	0.200	--	ND	0.869	--		1
4-Ethyltoluene	ND	0.200	--	ND	0.983	--		1
1,3,5-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
1,2,4-Trimethylbenzene	ND	0.200	--	ND	0.983	--		1
Benzyl chloride	ND	0.200	--	ND	1.04	--		1
1,3-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,4-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2-Dichlorobenzene	ND	0.200	--	ND	1.20	--		1
1,2,4-Trichlorobenzene	ND	0.200	--	ND	1.48	--		1
Naphthalene	ND	0.200	--	ND	1.05	--		1
Hexachlorobutadiene	ND	0.200	--	ND	2.13	--		1

## Lab Control Sample Analysis

### Batch Quality Control

**Project Name:** Q3 2024 SSDS MONITORING

**Project Number:** 01304

**Lab Number:** L2453073

**Report Date:** 09/24/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics in Air - Mansfield Lab Associated sample(s): 01-02 Batch: WG1974841-3								
Dichlorodifluoromethane	108		-		70-130	-		
Chloromethane	106		-		70-130	-		
Freon-114	117		-		70-130	-		
Vinyl chloride	110		-		70-130	-		
1,3-Butadiene	123		-		70-130	-		
Bromomethane	108		-		70-130	-		
Chloroethane	105		-		70-130	-		
Ethanol	110		-		40-160	-		
Vinyl bromide	96		-		70-130	-		
Acetone	101		-		40-160	-		
Trichlorofluoromethane	106		-		70-130	-		
Isopropanol	88		-		40-160	-		
1,1-Dichloroethene	109		-		70-130	-		
Tertiary butyl Alcohol	82		-		70-130	-		
Methylene chloride	107		-		70-130	-		
3-Chloropropene	108		-		70-130	-		
Carbon disulfide	107		-		70-130	-		
Freon-113	104		-		70-130	-		
trans-1,2-Dichloroethene	101		-		70-130	-		
1,1-Dichloroethane	104		-		70-130	-		
Methyl tert butyl ether	91		-		70-130	-		
2-Butanone	96		-		70-130	-		
cis-1,2-Dichloroethene	106		-		70-130	-		



## Lab Control Sample Analysis

### Batch Quality Control

**Project Name:** Q3 2024 SSDS MONITORING

**Project Number:** 01304

**Lab Number:** L2453073

**Report Date:** 09/24/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics in Air - Mansfield Lab Associated sample(s): 01-02 Batch: WG1974841-3								
Ethyl Acetate	110		-		70-130	-		
Chloroform	108		-		70-130	-		
Tetrahydrofuran	96		-		70-130	-		
1,2-Dichloroethane	104		-		70-130	-		
n-Hexane	114		-		70-130	-		
1,1,1-Trichloroethane	107		-		70-130	-		
Benzene	110		-		70-130	-		
Carbon tetrachloride	119		-		70-130	-		
Cyclohexane	114		-		70-130	-		
1,2-Dichloropropane	108		-		70-130	-		
Bromodichloromethane	119		-		70-130	-		
1,4-Dioxane	111		-		70-130	-		
Trichloroethene	107		-		70-130	-		
2,2,4-Trimethylpentane	114		-		70-130	-		
Heptane	108		-		70-130	-		
cis-1,3-Dichloropropene	115		-		70-130	-		
4-Methyl-2-pentanone	107		-		70-130	-		
trans-1,3-Dichloropropene	116		-		70-130	-		
1,1,2-Trichloroethane	108		-		70-130	-		
Toluene	92		-		70-130	-		
2-Hexanone	115		-		70-130	-		
Dibromochloromethane	117		-		70-130	-		
1,2-Dibromoethane	108		-		70-130	-		

## Lab Control Sample Analysis

### Batch Quality Control

**Project Name:** Q3 2024 SSDS MONITORING

**Project Number:** 01304

**Lab Number:** L2453073

**Report Date:** 09/24/24

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics in Air - Mansfield Lab Associated sample(s): 01-02 Batch: WG1974841-3								
Tetrachloroethene	100		-		70-130	-		
Chlorobenzene	106		-		70-130	-		
Ethylbenzene	100		-		70-130	-		
p/m-Xylene	102		-		70-130	-		
Bromoform	116		-		70-130	-		
Styrene	107		-		70-130	-		
1,1,2,2-Tetrachloroethane	113		-		70-130	-		
o-Xylene	104		-		70-130	-		
4-Ethyltoluene	104		-		70-130	-		
1,3,5-Trimethylbenzene	107		-		70-130	-		
1,2,4-Trimethylbenzene	108		-		70-130	-		
Benzyl chloride	102		-		70-130	-		
1,3-Dichlorobenzene	109		-		70-130	-		
1,4-Dichlorobenzene	110		-		70-130	-		
1,2-Dichlorobenzene	106		-		70-130	-		
1,2,4-Trichlorobenzene	105		-		70-130	-		
Naphthalene	94		-		70-130	-		
Hexachlorobutadiene	106		-		70-130	-		

**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

Serial\_No:09242417:00  
**Lab Number:** L2453073  
**Report Date:** 09/24/24

**Sample Receipt and Container Information**

Were project specific reporting limits specified? YES

**Cooler Information**

<b>Cooler</b>	<b>Custody Seal</b>
NA	Absent

**Container Information**

Container ID	Container Type	Cooler	Initial pH	Final pH	Temp deg C	Pres	Seal	Frozen Date/Time	Analysis(*)
L2453073-01A	Tedlar Bag 5 liter-Polypropylene Fitting	NA	NA			Y	Absent		TO15-LL(30)
L2453073-01X	Tedlar Bag 5 liter-Polypropylene Fitting	NA	NA			Y	Absent		TO15-LL(30)
L2453073-02A	Tedlar Bag 5 liter-Polypropylene Fitting	NA	NA			Y	Absent		TO15-LL(30)
L2453073-02X	Tedlar Bag 5 liter-Polypropylene Fitting	NA	NA			Y	Absent		TO15-LL(30)

\*Values in parentheses indicate holding time in days



**Project Name:** Q3 2024 SSDS MONITORING**Lab Number:** L2453073**Project Number:** 01304**Report Date:** 09/24/24

## 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: Data Usability Report*

**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

### 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.

**Chlordane:** 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.)

**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'.

**Gasoline Range Organics (GRO):** Gasoline Range Organics (GRO) results include all chromatographic peaks eluting from Methyl tert butyl ether through Naphthalene, with the exception of GRO analysis in support of State of Ohio programs, which includes all chromatographic peaks eluting from Hexane through Dodecane.

**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 for: PFHpA, PFHxS, PFOA, PFNA, PFDA and PFOS. For MassDEP DW compliance analysis only, the 'PFAS, Total (6)' result is defined as the summation of results at or above the RL. Note: If a 'Total' result is requested, the results of its individual components will also be reported.

**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. This represents an estimated concentration for Tentatively Identified Compounds (TICs).
- M** - Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.

**Report Format:** Data Usability Report



**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

**Data Qualifiers**

- ND** - Not detected at the reporting limit (RL) for the sample.
- 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.
- V** - The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- Z** - The batch matrix spike and/or duplicate associated with this target analyte has a recovery/RPD outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)

**Project Name:** Q3 2024 SSDS MONITORING  
**Project Number:** 01304

**Lab Number:** L2453073  
**Report Date:** 09/24/24

## REFERENCES

- 48 Compendium of Methods for the Determination of Toxic Organic Compounds in Ambient Air. Second Edition. EPA/625/R-96/010b, January 1999.

## 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 624.1:** m/p-xylene, o-xylene, Naphthalene

**EPA 625.1:** alpha-Terpineol

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

**EPA 8270E:** NPW: Dimethylnaphthalene, 1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; SCM: Dimethylnaphthalene, 1,4-Diphenylhydrazine.

**SM4500:** NPW: Amenable Cyanide; SCM: Total Phosphorus, TKN, NO<sub>2</sub>, NO<sub>3</sub>.

### Mansfield Facility

**SM 2540D:** TSS.

**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.

**Nonpotable Water:** **EPA RSK-175 Dissolved 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 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).

**Microbiology:** **SM9223B-Colilert-QT; Enterolert-QT, 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.



