

Date: May 1, 2025

- To: Mr. David G. Pratt, P.E. and Ms. Kathryn Lovell, New York State Department of Environmental Conservation, Region 8, East Avon-Lima Road, Avon, NY 14414
- From: Eric A. Warren, Roux Environmental Engineering and Geology, D.P.C.

#### Subject: March 2025 Monthly Progress Report Patriot Way Site No. 828223 293 Patriot Way, Chili, NY

Roux Environmental Engineering and Geology, D.P.C. (Roux) conducted a round of post injection well sampling on March 14, 2025. The sampling included monitoring wells MW-7B (located onsite north of the building) and MW-8 (located offsite to the west proximate to military base) which were sampled for target compound list (TCL) plus NYSDEC Commissioner Policy 51 (CP-51) VOCs. Please see the attached Alpha Analytical Report # L2516363 and the tabulated results comparing them to NYSDEC Division of Water Groundwater Quality Standards. We left the post remedial groundwater well sampling results in the chart for easy comparison and we marked the columns in green that are the most recent post injection results.

Please feel free to let me know if you have any questions.

Sincerely,

#### ROUX ENVIRONMENTAL ENGINEERING AND GEOLOGY, D.P.C.

Tric A. Warren

Eric A. Warren Senior Scientist II/Project Manager



TABLE 2

#### SUMMARY OF GROUNDWATER ANALYTICAL RESULTS PHASE II ENVIRONMENTAL INVESTIGATION 293 PATRIOT WAY ROCHESTER, NY

						Sample	Location								
PARAMETER <sup>1</sup>	GWQS <sup>2</sup>	MW-4B	MW-4B	MW-4B	MW-3	MW-3	MW-3	MW-B	MW-B	MW-B	TMW-1A	TMW-1A	TMW-1A	MW-7B	MW-8
		7/5/2024 <sup>3</sup>	10/24/2024 <sup>4</sup>	12/11/2024 <sup>5</sup>	7/5/2024 <sup>3</sup>	10/24/20244	12/11/2024 <sup>5</sup>	7/5/2024 <sup>3</sup>	10/24/2024 <sup>4</sup>	12/11/2024 <sup>5</sup>	7/5/2024 <sup>3</sup>	10/24/20244	12/11/2024 <sup>5</sup>	03/14/25	03/14/25
Volatile Organic Compounds	s (VOCs) - ug/L														
1,1-Dchloroethene	5	1.1	ND	ND	0.32 J	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.55
2-Butanone	50	ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	4.8 J	ND
Acetone	50	ND	ND	4.8 J	ND	ND	9.6	ND	ND	3.9 J	ND	1.5 J	ND	13	2.1 J
Benzene	1	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.16 J	ND	ND	ND	ND
Carbon disulfide		ND	ND	ND	ND	ND		ND	ND	ND	ND	ND	ND	1.9 J	ND
Cis-1,2-Dichloroethene	5	120	ND	2 J	100	ND	ND	1.7 J	1.2 J	ND	0.8 J	0.76 J	ND	ND	190
Cyclohexane		ND	ND	ND	ND	ND	ND	ND	ND	ND	ND	0.87 J	ND	ND	ND
Methyl Acetate		ND	ND	0.92 J	ND	ND	1.7 J	ND	ND	ND	ND	ND	ND	ND	ND
Methylcyclohexane		ND	ND	ND	ND	ND	ND	ND	1.4 J	ND	ND	1.5 J	ND	ND	ND
p-lsopropyltoluene	5	ND	ND	ND	ND	ND	ND	ND	ND	3.4	ND	ND	ND	ND	ND
Tetrachloroethene	5	ND	ND	ND	3.8	ND	ND	1.3	1.1	0.19 J	ND	ND	ND	ND	4.1
Trichloroethene	5	180	ND	8.8	82	ND	2.2	11	9.4	1.4	3.1	4.2	4.3	ND	160 D
Trans-1,2- Dichloroethene	5	25	ND	ND	2.6	ND	ND	ND	ND	ND	ND	ND	ND	ND	4
Vinyl Chloride	2	10	ND	0.47 J	0.66 J	ND	ND	ND	ND	ND	ND	ND	ND	ND	1.4
Dissolved Metals - ug/L															
Iron Total		0.000757	0.00711	4.8	0.000224	0.0178	3.29	0.108	0.118	23.6	0.0976	0.0494	13.1		
Iron Dissolved		0.0002	0.0025	0.521	0.0000315 J	0.00126 J	ND	0.0000574	0.00502	0.0215 J	0.0000646	0.000758	ND		
General Chemistry - ug/L															
Nitrogen, Nitrate/ Nitrite		ND	ND	ND	0.00019	ND	ND	0.0017	0.00138	0.309	0.013	0.0213	13.5		
Total Organic Carbon		0.00195	0.15	47	0.00248	0.36	94	0.00532	0.0039	2.4	0.0263	0.011	7.4		
Anions- ug/L															
Sulfate		0.0677	0.032	6.5 J	0.0428	0.031	1.7 J	0.0284	0.022	29	0.0347	0.031	28		

Notes:
1. Only those parameters detected at a minimum of one sample location are presented in this table; all other compounds were reported as non-detect.
2. Values per NYSDEC Division of Water Ambient Water Quality Standards and Guidance Values and Groundwater Effluent Limitations - Class GA (TOGS 1.1.1)
3. Total reported as a constraint instance of the same constraint of the same constraint instance of the sa

4. Post Injection Analytical Report L2462191
 5. Post Injection Analytical Report L2472614
 5. Post Injection Analytical Report L2576363
 Definitions:
 ND = Parameter not detected above laboratory detection limit.
 \*\_\* = Sample not analyzed for parameter or no SCO available for the parameter.
 J = Estimated Value - Below calibration range.
 D = Concentration of analyte was quantified from diluted analysis.
 BOLD = Result exceeds GWQS.



### ANALYTICAL REPORT

Lab Number:	L2516363
Client:	Roux
	2558 Hamburg Turnpike
	Suite 300
	Buffalo, NY 14218
ATTN:	Eric Warren
Phone:	(716) 856-0599
Project Name:	KADDIS
Project Number:	4351.0001B000
Report Date:	03/28/25

The original project report/data package is held by Pace Analytical Services. This report/data package is paginated and should be reproduced only in its entirety. Pace Analytical Services 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).

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

Serial\_No:03282510:00

Project Name:KADDISProject Number:4351.0001B000

 Lab Number:
 L2516363

 Report Date:
 03/28/25

Lab Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2516363-01	MW-7B	WATER	293 PATRIOT WAY	03/14/25 18:00	03/20/25
L2516363-02	MW-8	WATER	293 PATRIOT WAY	03/14/25 17:00	03/20/25



Project Name: KADDIS Project Number: 4351.0001B000 
 Lab Number:
 L2516363

 Report Date:
 03/28/25

#### **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 Pace 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, Pace'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 Pace Project Manager and made arrangements for Pace 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: KADDIS Project Number: 4351.0001B000 
 Lab Number:
 L2516363

 Report Date:
 03/28/25

#### **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.

Volatile Organics

L2516363-02: The sample was received in the proper acid-preserved containers. Upon analysis the pH was determined to be greater than 2 and the method required holding time was exceeded.

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.

Melissa Sturgis Melissa Sturgis

Authorized Signature:

Title: Technical Director/Representative

Date: 03/28/25

# ORGANICS



## VOLATILES



			Serial_No:03282510:00			
Project Name:	KADDIS		Lab Number:	L2516363		
Project Number:	4351.0001B000		Report Date:	03/28/25		
		SAMPLE RESULTS				
Lab ID:	L2516363-01		Date Collected:	03/14/25 18:00		
Client ID:	MW-7B		Date Received:	03/20/25		
Sample Location:	293 PATRIOT WAY		Field Prep:	Not Specified		
Sample Depth:						
Matrix:	Water					
Analytical Method:	1,8260D					
Analytical Date:	03/26/25 00:08					
Analyst:	MKS					

Parameter	Result	Qualifier	Units	RL	MDL	<b>Dilution Factor</b>
Volatile Organics by GC/MS - Wes	tborough 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

						Serial_No	0:03282510:00
Project Name:	KADDIS				Lab Nu	ımber:	L2516363
Project Number:	4351.0001B000				Report	Date:	03/28/25
-		SAMP	LE RESULTS	6			
Lab ID: Client ID: Sample Location:	L2516363-01 MW-7B 293 PATRIOT WAY			Date Collected: Date Received: Field Prep:		03/14/25 18:00 03/20/25 Not Specified	
Sample Depth:							
Parameter		Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics b	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
Dichlorodifluoromethan	e	ND		ug/l	5.0	1.0	1
Acetone		13		ug/l	5.0	1.5	1
Carbon disulfide		1.9	J	ug/l	5.0	1.0	1
2-Butanone		4.8	J	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-chloropi	ropane	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	9	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	Acceptance Qualifier Criteria
1,2-Dichloroethane-d4	107	70-130
Toluene-d8	100	70-130
4-Bromofluorobenzene	97	70-130
Dibromofluoromethane	115	70-130

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			Serial_No:03282510:00		
Project Name:	KADDIS		Lab Number:	L2516363	
Project Number:	4351.0001B000		Report Date:	03/28/25	
		SAMPLE RESULTS			
Lab ID: Client ID: Sample Location:	L2516363-02 MW-8 293 PATRIOT WAY		Date Collected: Date Received: Field Prep:	03/14/25 17:00 03/20/25 Not Specified	
Sample Depth: Matrix: Analytical Method: Analytical Date: Analyst:	Water 1,8260D 03/26/25 00:32 MKS				

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	4.1		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	1.4		ug/l	1.0	0.07	1		
Chloroethane	ND		ug/l	2.5	0.70	1		
1,1-Dichloroethene	0.55		ug/l	0.50	0.17	1		
trans-1,2-Dichloroethene	4.0		ug/l	2.5	0.70	1		
Trichloroethene	220	Е	ug/l	0.50	0.18	1		
1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1		



Serial_No:03282510:						0:03282510:00	
Project Name:	KADDIS				Lab Nu	umber:	L2516363
Project Number:	4351.0001B000				Report	Date:	03/28/25
-		SAMP	LE RESULTS	5			
Lab ID: Client ID: Sample Location:	L2516363-02 MW-8 293 PATRIOT WAY				Date Co Date Re Field Pre	ceived:	03/14/25 17:00 03/20/25 Not Specified
Sample Depth:							
Parameter		Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics b	y GC/MS - Westborough	Lab					
1.2 Disblarchanzana		ND			2.5	0.70	4
1,3-Dichlorobenzene		ND		ug/l			1
1,4-Dichlorobenzene				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		190		ug/l	2.5	0.70	1
Styrene		ND		ug/l	2.5	0.70	1
Dichlorodifluoromethane	9	ND	•	ug/l	5.0	1.0	1
Acetone		2.1	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-chloropro	opane	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	106		70-130	
Toluene-d8	100		70-130	
4-Bromofluorobenzene	95		70-130	
Dibromofluoromethane	109		70-130	

				Serial_No	0:03282510:00
Project Name:	KADDIS			Lab Number:	L2516363
Project Number:	4351.0001B000			Report Date:	03/28/25
			SAMPLE RESULTS		
Lab ID:	L2516363-02	D		Date Collected:	03/14/25 17:00
Client ID:	MW-8			Date Received:	03/20/25
Sample Location:	293 PATRIOT WAY			Field Prep:	Not Specified
Sample Depth:					
Matrix:	Water				
Analytical Method:	1,8260D				
Analytical Date:	03/27/25 13:13				
Analyst:	MKS				

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
olatile Organics by GC/MS - Westboroug	h Lab					
richloroethene	160		ug/l	2.0	0.70	4
Surrogate			% Recovery	Qualifier		eptance riteria
1,2-Dichloroethane-d4			103			70-130
Toluene-d8			97			70-130
4-Bromofluorobenzene			102			70-130
Dibromofluoromethane			113			70-130

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Project Name:KADDISProject Number:4351.0001B000

 Lab Number:
 L2516363

 Report Date:
 03/28/25

## Method Blank Analysis Batch Quality Control

Analytical Method:1,8260Analytical Date:03/25/Analyst:TMS

1,8260D 03/25/25 23:06 TMS

arameter	Result	Qualifier Units	RL	MDL
olatile Organics by GC/MS - W	/estborough Lab	for sample(s): 01-02	Batch:	WG2045483-5
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

Pace

Project Name: KADDIS Project Number: 4351.0001B000

Lab Number: L2516363 **Report Date:** 03/28/25

## Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260D Analytical Date: Analyst: TMS

03/25/25 23:06

arameter	Result C	Qualifier Units	RL	MDL
olatile Organics by GC/MS -	Westborough Lab fo	or sample(s): 01-02	Batch:	WG2045483-5
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

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 Lab Number:
 L2516363

 Report Date:
 03/28/25

Project Name:KADDISProject Number:4351.0001B000

## Method Blank Analysis Batch Quality Control

Analytical Method:1,8260DAnalytical Date:03/25/25 23:06Analyst:TMS

Parameter	Result	Qualifier	Units	5	RL	MDL	
Volatile Organics by GC/MS	- Westborough Lab	o for sample	e(s):	01-02	Batch:	WG2045483-5	

Surrogate	%Recovery	Qualifier	Criteria		
4.0 Dishlarathara da	110		70.400		
1,2-Dichloroethane-d4	110		70-130		
Toluene-d8	100		70-130		
4-Bromofluorobenzene	96		70-130		
Dibromofluoromethane	112		70-130		

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Project Name:KADDISProject Number:4351.0001B000

 Lab Number:
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 Report Date:
 03/28/25

## Method Blank Analysis Batch Quality Control

Analytical Method:1,8260DAnalytical Date:03/27/25 12:51Analyst:PID

arameter	Result	Qualifier Units	s RL	MDL	
olatile Organics by GC/MS -	Westborough Lab	for sample(s):	02 Batch:	WG2046075-5	
1,1-Dichloroethane	ND	ug/l	2.5	0.70	
Tetrachloroethene	ND	ug/l	0.50	0.18	
1,2-Dichloroethane	ND	ug/l	0.50	0.13	
1,1,1-Trichloroethane	ND	ug/l	2.5	0.70	
Benzene	ND	ug/l	0.50	0.16	
Toluene	ND	ug/l	2.5	0.70	
Ethylbenzene	ND	ug/l	2.5	0.70	
Vinyl chloride	ND	ug/l	1.0	0.07	
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	
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	
1,2-Dibromoethane	ND	ug/l	2.0	0.65	
Isopropylbenzene	ND	ug/l	2.5	0.70	

		Acceptance		
Surrogate	%Recovery	Qualifier	Criteria	
1,2-Dichloroethane-d4	108		70-130	
Toluene-d8	95		70-130	
4-Bromofluorobenzene	99		70-130	
Dibromofluoromethane	117		70-130	

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Project Name:KADDISProject Number:4351.0001B000

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**Report Date:** 03/28/25

arameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
olatile Organics by GC/MS - Westborou	gh Lab Associa	ted sample(s):	01-02 Bate	ch: WG204	5483-3 WG204	5483-4		
Methylene chloride	88		89		70-130	1		20
1,1-Dichloroethane	100		100		70-130	0		20
Chloroform	98		99		70-130	1		20
Carbon tetrachloride	100		99		63-132	1		20
1,2-Dichloropropane	97		95		70-130	2		20
Dibromochloromethane	89		88		63-130	1		20
1,1,2-Trichloroethane	85		86		70-130	1		20
Tetrachloroethene	98		96		70-130	2		20
Chlorobenzene	100		99		75-130	1		20
Trichlorofluoromethane	91		92		62-150	1		20
1,2-Dichloroethane	95		96		70-130	1		20
1,1,1-Trichloroethane	100		100		67-130	0		20
Bromodichloromethane	100		100		67-130	0		20
trans-1,3-Dichloropropene	82		84		70-130	2		20
cis-1,3-Dichloropropene	98		98		70-130	0		20
Bromoform	76		77		54-136	1		20
1,1,2,2-Tetrachloroethane	80		83		67-130	4		20
Benzene	89		89		70-130	0		20
Toluene	96		94		70-130	2		20
Ethylbenzene	100		100		70-130	0		20
Chloromethane	82		81		64-130	1		20
Bromomethane	38	Q	38	Q	39-139	0		20
Vinyl chloride	83		85		55-140	2		20



Project Name:KADDISProject Number:4351.0001B000

Lab Number: L2516363

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arameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits
olatile Organics by GC/MS - Westbord	ough Lab Associat	ed sample(s)	: 01-02 Batcl	h: WG204	45483-3 WG204	5483-4		
Chloroethane	76		73		55-138	4		20
1,1-Dichloroethene	88		90		61-145	2		20
trans-1,2-Dichloroethene	95		94		70-130	1		20
Trichloroethene	95		95		70-130	0		20
1,2-Dichlorobenzene	96		96		70-130	0		20
1,3-Dichlorobenzene	100		100		70-130	0		20
1,4-Dichlorobenzene	100		98		70-130	2		20
Methyl tert butyl ether	83		84		63-130	1		20
p/m-Xylene	105		105		70-130	0		20
o-Xylene	100		105		70-130	5		20
cis-1,2-Dichloroethene	96		98		70-130	2		20
Styrene	100		100		70-130	0		20
Dichlorodifluoromethane	68		68		36-147	0		20
Acetone	93		90		58-148	3		20
Carbon disulfide	94		94		51-130	0		20
2-Butanone	90		94		63-138	4		20
4-Methyl-2-pentanone	77		70		59-130	10		20
2-Hexanone	86		89		57-130	3		20
Bromochloromethane	94		96		70-130	2		20
1,2-Dibromoethane	84		86		70-130	2		20
1,2-Dibromo-3-chloropropane	82		85		41-144	4		20
Isopropylbenzene	110		110		70-130	0		20
1,2,3-Trichlorobenzene	95		97		70-130	2		20



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LCSD LCS %Recovery RPD %Recovery %Recovery Limits Limits Parameter Qual Qual RPD Qual Volatile Organics by GC/MS - Westborough Lab Associated sample(s): 01-02 Batch: WG2045483-3 WG2045483-4 1,2,4-Trichlorobenzene 96 98 70-130 2 20 Methyl Acetate 97 87 70-130 20 11 Cyclohexane 110 110 70-130 0 20 1,4-Dioxane 92 96 56-162 20 4 Freon-113 95 96 70-130 20 1 Methyl cyclohexane 70-130 20 110 110 0

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	99	103	70-130
Toluene-d8	97	98	70-130
4-Bromofluorobenzene	96	98	70-130
Dibromofluoromethane	101	104	70-130



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Parameter	LCS %Recovery	Qual	LCSD %Recovery	%Recovery Qual Limits	RPD	RPD Qual Limits	
/olatile Organics by GC/MS - Westborou	gh Lab Associate	ed sample(s):	02 Batch:	WG2046075-3 WG2046	075-4		
1,1-Dichloroethane	100		98	70-130	2	20	
Tetrachloroethene	100		100	70-130	0	20	
1,2-Dichloroethane	100		100	70-130	0	20	
1,1,1-Trichloroethane	100		100	67-130	0	20	
Benzene	90		88	70-130	2	20	
Toluene	100		98	70-130	2	20	
Ethylbenzene	93		92	70-130	1	20	
Vinyl chloride	77		79	55-140	3	20	
1,1-Dichloroethene	98		90	61-145	9	20	
trans-1,2-Dichloroethene	100		99	70-130	1	20	
Trichloroethene	99		99	70-130	0	20	
Methyl tert butyl ether	67		70	63-130	4	20	
p/m-Xylene	100		95	70-130	5	20	
o-Xylene	100		95	70-130	5	20	
cis-1,2-Dichloroethene	100		96	70-130	4	20	
1,2-Dibromoethane	110		110	70-130	0	20	
Isopropylbenzene	93		87	70-130	7	20	



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Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	Qual	RPD Limits	
Volatile Organics by GC/MS - Westboroug	gh Lab Associat	ed sample(s)	: 02 Batch:	WG204607	75-3 WG204607	5-4			

Surrogate	LCS %Recovery Qual	LCSD %Recovery Qual	Acceptance Criteria
1,2-Dichloroethane-d4	102	103	70-130
Toluene-d8	103	102	70-130
4-Bromofluorobenzene	97	96	70-130
Dibromofluoromethane	104	102	70-130



#### Project Name: KADDIS *Project Number:* 4351.0001B000

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#### Sample Receipt and Container Information

Were project specific reporting limits specified?

YES

#### **Cooler Information**

Cooler	Custody Seal						
A	Absent						

#### Container Information

Container Information		Initial	Final Temp				Frozen			
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)	
L2516363-01A	Vial HCI preserved	А	NA		2.4	Y	Absent		NYTCL-8260-R2(14)	
L2516363-01B	Vial HCI preserved	А	NA		2.4	Υ	Absent		NYTCL-8260-R2(14)	
L2516363-01C	Vial HCI preserved	А	NA		2.4	Y	Absent		NYTCL-8260-R2(14)	
L2516363-02A	Vial HCI preserved	А	NA		2.4	Y	Absent		NYTCL-8260-R2(14)	
L2516363-02B	Vial HCI preserved	А	NA		2.4	Y	Absent		NYTCL-8260-R2(14)	
L2516363-02C	Vial HCI preserved	А	NA		2.4	Y	Absent		NYTCL-8260-R2(14)	



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## Project Name: KADDIS

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#### GLOSSARY

#### Acronyms

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	<ul> <li>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.)</li> </ul>
	Limit of Quantitation: The value at which an instrument can accurately measure an analyte at a specific concentration. The LOQ includes any adjustments from dilutions, concentrations or moisture content, where applicable. (DoD report formats only.)
MDL	- Method Detection Limit: This value represents the level to which target analyte concentrations are reported as estimated values, when those target analyte concentrations are quantified below the reporting limit (RL). The MDL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
MS	- Matrix Spike Sample: A sample prepared by adding a known mass of target analyte to a specified amount of matrix sample for which an independent estimate of target analyte concentration is available. For Method 332.0, the spike recovery is calculated using the native concentration, including estimated values.
MSD	- Matrix Spike Sample Duplicate: Refer to MS.
NA	- Not Applicable.
NC	- Not Calculated: Term is utilized when one or more of the results utilized in the calculation are non-detect at the parameter's reporting unit.
NDPA/DPA	- N-Nitrosodiphenylamine/Diphenylamine.
NI	- Not Ignitable.
NP	- Non-Plastic: Term is utilized for the analysis of Atterberg Limits in soil.
NR	- No Results: Term is utilized when 'No Target Compounds Requested' is reported for the analysis of Volatile or Semivolatile Organic TIC only requests.
RL	- Reporting Limit: The value at which an instrument can accurately measure an analyte at a specific concentration. The RL includes any adjustments from dilutions, concentrations or moisture content, where applicable.
RPD	- Relative Percent Difference: The results from matrix and/or matrix spike duplicates are primarily designed to assess the precision of analytical results in a given matrix and are expressed as relative percent difference (RPD). Values which are less than five times the reporting limit for any individual parameter are evaluated by utilizing the absolute difference between the values; although the RPD value will be provided in the report.
SRM	- Standard Reference Material: A reference sample of a known or certified value that is of the same or similar matrix as the associated field samples.
STLP	- Semi-dynamic Tank Leaching Procedure per EPA Method 1315.
TEF	- Toxic Equivalency Factors: The values assigned to each dioxin and furan to evaluate their toxicity relative to 2,3,7,8-TCDD.
TEQ	- Toxic Equivalent: The measure of a sample's toxicity derived by multiplying each dioxin and furan by its corresponding TEF and then summing the resulting values.
TIC	- Tentatively Identified Compound: A compound that has been identified to be present and is not part of the target compound list (TCL) for the method and/or program. All TICs are qualitatively identified and reported as estimated concentrations.

Report Format: DU Report with 'J' Qualifiers



## Project Name: KADDIS

#### **Project Number:** 4351.0001B000

### Lab Number: L2516363 Report Date: 03/28/25

#### Footnotes

•	-	•	•••	•	-

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

#### Terms

1

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 Waterpreserved 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(a)+(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, 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 concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the concentrations of the analyte at less than ten times (10x) the 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 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

Pace Analytical Services 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 Pace Analytical Services shall be to re-perform the work at it's own expense. In no event shall Pace Analytical Services 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 Pace Analytical Services.

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 – 8 Walkup Dr. Westborough, MA 01581

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

EPA 625.1: alpha-Terpineol

**EPA 8260D:** <u>NPW</u>: 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene; <u>SCM</u>: lodomethane (methyl iodide), 1,2,4,5-Tetramethylbenzene; 4-Ethyltoluene. **EPA 8270E:** <u>NPW</u>: Dimethylnaphthalene,1,4-Diphenylhydrazine, alpha-Terpineol, Azobenzene; <u>SCM</u>: Dimethylnaphthalene,1,4-Diphenylhydrazine. **SM4500**: <u>NPW</u>: Amenable Cyanide; <u>SCM</u>: Total Phosphorus, TKN, NO2, NO3.

#### Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048

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. MADEP-APH. Nonpotable Water: EPA RSK-175 Dissolved Gases

Biological Tissue Matrix: EPA 3050B

Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048 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

#### The following test method is not included in our New Jersey Secondary NELAP Scope of Accreditation:

Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048 Determination of Selected Perfluorinated Alkyl Substances by Solid Phase Extraction and Liquid Chromatography/Tandem Mass Spectrometry Isotope Dilution (via Alpha SOP 23528)

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

#### Westborough Facility - 8 Walkup Dr. Westborough, MA 01581

**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 – 320 Forbes Blvd. Mansfield, MA 02048

#### **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** 

#### **Certification IDs:**

#### Westborough Facility – 8 Walkup Dr. Westborough, MA 01581

CT PH-0826, IL 200077, IN C-MA-03, KY JY98045, ME MA00086, MD 348, MA M-MA086, NH 2064, NJ MA935, NY 11148, NC (DW) 25700, NC (NPW/SCM) 666, OR MA-1316, PA 68-03671, RI LAO00065, TX T104704476, VT VT-0935, VA 460195

#### Mansfield Facility – 320 Forbes Blvd. Mansfield, MA 02048

CT PH-0825, ANÅB/DoD L2474, IL 200081, IN C-MA-04, KY KY98046, LA 3090, ME MA00030, MI 9110, MN 025-999-495, NH 2062, NJ MA015, NY 11627, NC (NPW/SCM) 685, OR MA-0262, PA 68-02089, RI LAO00299, TX T-104704419, VT VT-0015, VA 460194, WA C954

#### Mansfield Facility – 120 Forbes Blvd. Mansfield, MA 02048

ANAB/DoD L2474, ME MA01156, MN 025-999-498, NH 2249, NJ MA025, NY 12191, OR 4203, TX T104704583, VA 460311, WA C1104.

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

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