

February 1, 2023

Mr. Charles T. Gregory, LG Engineering Geologist 1 Section C, Bureau E Division of Environmental Remediation 625 Broadway Albany, NY 12233-7011

Re: Alco Manufacturing Corporation, LLC

NYSDEC Site No. 826011

September 2022 Post-Remedial Groundwater Monitoring Report-REVISED February 2023

Dear Mr. Gregory:

Benchmark Civil/Environmental Engineering & Geology, PLLC (Benchmark), has prepared this letter report to transmit the results of the September 2022 (Revised February 2023) post-remedial groundwater monitoring event at the former Kaddis Manufacturing (Kaddis) Site located in Lima, NY (see Figure 1). In November 2021, Alco Manufacturing Corporation, LLC (Alco) acquired the Site from Kaddis. The Site is currently owned and operated by Alco.

FIELD SAMPLING PROCEDURE

In accordance with Benchmark's revised October 3, 2008 Work Plan, four groundwater monitoring locations were designated for sampling during the subject 15-month sampling event: MW-3, MW-5, MW-201D, and the former Supply Well. Benchmark has continued to use passive diffusion bags (PDBs). The PDB sampler is a semi-permeable, low-density polyethylene membrane designed to allow volatile organic compounds (VOCs) to flow into the PDB until equilibrium is reached between the formation and the PDB.

Benchmark field staff deployed the PDBs on August 18, 2022 and retrieved and sampled the PDBs on September 1, 2022. Attachment 1 includes PDB deployment and retrieval logs. The groundwater samples were transferred to laboratory supplied, pre-preserved sample vials and transported, under chain of custody control, to Alpha Analytical (Alpha) in Westborough, MA for analysis of Target Compound List (TCL) VOCs per USEPA Method 8260B.

ANALYTICAL RESULTS

Attachment 2 includes the analytical data package from Alpha. Table 1 summarizes the detected compounds and compares the results to NY State Groundwater Quality Standards and Guidance Values (NYSDEC TOGS 1.1.1, June 1998). As indicated on Table 1, VOC detections were generally limited to trace concentrations (below 1 part per million) except for trichloroethene (TCE) and (CIS-1,2-Dichloroethene) at well MW-201D.

GROUNDWATER FLOW DIRECTION

On September 1, 2022, groundwater levels were measured in all on-site wells (MW-1 through MW-6, MW-201D, MW-202, and the Supply Well). Groundwater elevations are summarized on Table 2 and

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presented as an isopotential map on Figure 1. The Supply Well, with a total depth of 185 feet below ground surface (fbgs), reflects the deeper groundwater aquifer, and was therefore not used to develop the isopotential map. In addition, an artificial mound has been historically observed at well MW-201D and is thought to be caused from unconsolidated structural fill materials that were used to backfill the remedial excavation in this area of the Site. As such, the water level collected from well MW-201D is reflective of the trapped water. As shown on Figure 1, groundwater flow is generally to the north and northwest, with a localized component flowing northeast toward Honeoye Creek on the east side of the Site consistent with previous studies.

HISTORICAL COMPARISONS

Attachment 3 graphically depicts the September 2022 total VOC concentrations at each of the sampled locations with historical concentrations for key parameters including 1,1,1-trichloroethane (1,1,1-TCA), 1,1-dichloroethene (1,1-DCE), cis-1,2-dichloroethene (cis-1,2-DCE), tetrachloroethene (PCE), and TCE.

The September 2022 data indicates a slight increase in the total concentration of VOCs at wells MW-5 and MW-201D when compared to May 2021 sampling event, however concentrations are consistent with the historic fluctuating seasonal trends. Conversely, the total VOC concentrations detected at MW-3 and the Supply Well during the September 2022 event slightly decreased as compared to the May 2021 sampling event. These fluctuations are again consistent with historical sampling trends. Overall the concentrations detected at all of the sampling locations are well below historic highs and indicate that natural attenuation processes continue to reduce downgradient concentrations and mitigate associated off-site environmental impact.

The electronic data delivery (EDD) format was uploaded to NYSDEC's EQuIS database on October 18th, 2022. The next sampling event is expected to take place in November/December of 2023.

MONITORING WELL REPAIRS

On November 4, 2022, Benchmark's subcontract driller, Trec Environmental, Inc. repaired damaged road box covers at monitoring wells MW-202 and MW-201D. Photo-documentation of the repairs is provided as Attachment 4.

Please contact us with any questions.

Sincerely,

Benchmark Civil/Environmental Engineering & Geology, PLLC

Thomas H. Forbes, P.E.

President

Att.

ec: Robert Papenfuss (Alco)

T. Behrendt (Benchmark)



TABLES





TABLE 1

POST-REMEDIAL GROUNDWATER MONITORING RESULTS September 2022

Enarc-O Machine Products, Inc. Lima, New York NYSDEC Registry No. 8-26-011

PARAMETER ¹	MW-3	MW-201D	MW-5	SUPPLY WELL	GWQS ²			
Volatile Organic Compounds (ug/L):								
1,1,1-Trichloroethane	8.1	160	4 J	ND	5			
1,1-Dichloroethane	ND	ND	3.4 J	ND	5			
1,1-Dichloroethene	3.4	70	2.4	0.27 J	5			
cis-1,2-Dichloroethene	24	1200	140	1.1 J	5			
Tetrachloroethene	5.3	54	5.3	ND	5			
Trichloroethene	290	6,000	500	12	5			
Total VOCs	330.8	7484	655.1	13.37				

Notes:

- 1. Only those compounds detected above the method detection limit at a minimum of one sample location are reported in this table.
- 2. NYSDEC Class "GA" Groundwater Quality Standards (GWQS) as per 6 NYCRR Part 703. Guidance value used when Standard value not available.

Acronyms:

ND = Parameter was not detected above laboratory reporting limit.

J = Indicates an estimated value.

NA = Not Analyzed.

BOLD

= Value exceeds GWQS.



TABLE 2

SUMMARY OF GROUNDWATER ELEVATIONS September 1, 2022

Enarc-O Machine Products, Inc. Lima, New York NYSDEC Registry No. 8-26-011

WELL ID	Depth to Water	TOR Elevation ¹	Groundwater Elevation	Bottom Depth
MW - 1	26.51	502.69	476.18	35.83
MW - 2	27.74	506.79	479.05	33.78
MW - 3	26.28	505.27	478.99	34.58
MW - 4	23.73	500.73	477	34.32
MW - 5	23.70	502.38	478.68	30.49
MW - 6	27.72	504.86	477.14	37.88
MW - 202	25.03	504.28	479.25	34.81
MW - 201D	14.24	501.04	486.8	29.35
Supply Well	109.00	503.39	394.39	185.00

Notes:

1. Top of riser survey was completed on 5/31/2016.

FIGURE



ATTACHMENT 1

PDB Logs





GROUNDWATER WELL PDB COLLECTION & RECOVERY LOG

(PASSIVE DIFFUSION BAG)

Project Name:	Kaddis Enarco	WELL NUMBER: MW-3
Project Number	: 0127-021-101	Sample Matrix: Water
Client:	Kaddis, Mfg	Weather: Denny 76'S
WELL DAT	TA:	
Casing Diame	eter (inches): 4.0	Casing Material: PVC
Screened inte	erval (fbTOR):	Screen Material:
Static Water I	Level (fbTOR): 29,14	Bottom Depth (fbTOR): 34.57
Elevation Top	of Well Riser (fmsl): 505.27	Ground Surface Elevation (fmsl): 503.40
Elevation Top	of Screen (fmsl):	Stick-up (feet): 1.87'
PDB DATA	l:	
	in well (fbTOR):	Is PDB harness and line dedicated to sample location? yes no
Condition of V	Vell: Gove	Is PDB located at aproximatley center of screen or water column?
Field Personn	Vell: Good nel: TA3	
Installatio	on.	
Date of PDB		5122
Time of PDB	111	12
Retrieval		
Date of PDB r	retrieval. 9/11	122 WL 26.28
Time of PDB	retrieval: UL8	
Condition of F	PDB:	
f PDR conta	iins visible sediment, check l	PDR integrity and re-sample
. , DB comu	and violate deathfort, effect t	bb mognly and re sample.
COMMENT	S :	

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GROUNDWATER WELL PDB COLLECTION & RECOVERY LOG

(PASSIVE DIFFUSION BAG)

Kaddis Enarco WELL NUMBER: **MW-5** Project Name: 0127-021-101 Sample Matrix: Water Project Number: Weather: Client: Kaddis Mfg **WELL DATA:** 4.0 **PVC** Casing Diameter (inches): Casing Material: Screened interval (fbTOR): hole Screen Material: Open 26.35, 27.63 _30.03 Bottom Depth (fbTOR): Static Water Level (fbTOR): Elevation Top of Well Riser (fmsl): 502.38 Ground Surface Elevation (fmsl): Elevation Top of Screen (fmsl): Stick-up (feet): 1.9 PDB DATA: Is PDB harness and line dedicated to sample location? Depth of PDB in well (fbTOR): Condition of Well: See Note Is PDB located at aproximatley center of screen or water column? Field Personnel: THIS Installation: 8/18/22 Date of PDB placement. Time of PDB placement: 1200 Retrieval: WL -23.70 11/22 Date of PDB retrieval. Time of PDB retrieval: Condition of PDB: If PDB contains visible sediment, check PDB integrity and re-sample. 6f Sediment to Eacility PDB Placement 1.3-

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GROUNDWATER WELL PDB COLLECTION & RECOVERY LOG

(PASSIVE DIFFUSION BAG)

Project Name:	Kaddis Enarco		WELL NUMBER:	MW-201D
Project Number:	0127-001-104		Sample Matrix:	Water
Client:	Kaddis, Mfg		Weather: 500	my Mid 76"
WELL DATA	4 :			
Casing Diamete	er (inches): 2.0)	Casing Material:	PVC
Screened inter	val (fbTOR):		Screen Material:	open hole
Static Water Le	evel (fbTOR):	1.20	Bottom Depth (fbTOR	
Elevation Top of	of Well Riser (fmsl):	501.44	Ground Surface Eleva	ation (fmsl); 501.33
Elevation Top of	of Screen (fmsl):	14	Stick-up (feet):	Flush
PDB DATA:				
Depth of PDB is			Is PDB harness and li	ine dedicated to sample location?
Condition of W	ell: 9002		Is PDB located at apr	oximatley center of screen or water column?
Field Personne				
Installation		8/14/2	7.	
Time of PDB pl		8/18/2		
		1104		
Retrieval:				
Date of PDB re	trieval.	9/1/2	2	14.24
Time of PDB re	trieval:	1143		
Condition of PE	DB:	500		
f PDB contail	ns visible sedin	-	DB integrity and re-	sample.
		·	•	
COMMENTS	S:			

PREPARED BY:



GROUNDWATER WELL PDB COLLECTION & RECOVERY LOG

(PASSIVE DIFFUSION BAG)

Project Name:	Kaddis Enarco	WELL NUMBER:	Supply Well
Project Number:	0127-001-104	Sample Matrix:	Water
Client:	Kaddis, Mfg	Weather:	Survey 70°5
NELL DATA	4 :		A.
Casing Diamet	er (inches): 6.0	Casing Material: Ste	el
Screened inter	val (fbTOR):	Screen Material:	pen hole
Static Water Le	evel (fbTOR): 107.72	Bottom Depth (fbTOR):	185.00 53
Elevation Top	of Well Riser (fmsl): 503.39	Ground Surface Elevation	
Elevation Top	of Screen (fmsl):	Stick-up (feet): 1.2	
DDD DATA			
PDB DATA:	n well (fbTOR):	Is PDB harness and line	dedicated to sample location? yes no
Condition of W			matley center of screen or water column?
Field Personne	5002	13 1 DD located at aproxi	~ No /
Ticla i ersonire	11 743		780
Installatio	n:		
Date of PDB pl	acement. 8/18	122	
Time of PDB pl		70	
Retrieval:	1		.
Date of PDB re	trieval.	22	109.8
	trieval: U27		
Time of PDB re	NZr		

PREPARED BY:

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ATTACHMENT 2

ALPHA ANALYTICAL LABORATORY DATA PACKAGE

SEPTEMBER 2022





ANALYTICAL REPORT

Lab Number: L2247717

Client: Benchmark & Turnkey Companies

2558 Hamburg Turnpike

Suite 300

Buffalo, NY 14218

ATTN: Thomas Behrendt Phone: (716) 856-0599

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Report Date: 09/16/22

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

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

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



Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Lab Number: L2247717 **Report Date:** 09/16/22

Alpha Sample ID	Client ID	Matrix	Sample Location	Collection Date/Time	Receive Date
L2247717-01	MW-3	WATER	HONEOYE FALLS, NY	09/01/22 11:18	09/01/22
L2247717-02	MW-5	WATER	HONEOYE FALLS, NY	09/01/22 11:06	09/01/22
L2247717-03	MW-201D	WATER	HONEOYE FALLS, NY	09/01/22 11:43	09/01/22
L2247717-04	SUPPLY WELL	WATER	HONEOYE FALLS, NY	09/01/22 11:22	09/01/22
L2247717-05	TRIP BLANK	WATER	HONEOYE FALLS, NY	09/01/22 00:00	09/01/22



L2247717

Lab Number:

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001 **Report Date:** 09/16/22

Case Narrative

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

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

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

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

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

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



Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

Case Narrative (continued)

Report Submission

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

Sample Receipt

L2247717-04: The collection date and time on the chain of custody was 01-SEP-22 11:22; however, the collection date/time on the container label was 01-SEP-22 11:27. At the client's request, the collection date/time is reported as 01-SEP-22 11:22.

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:

Title: Technical Director/Representative Date: 09/16/22

Sufani Morrissey-Tiffani Morrissey

ORGANICS



VOLATILES



09/01/22 11:18

Not Specified

09/01/22

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

SAMPLE RESULTS

Lab Number: L2247717

Report Date: 09/16/22

Date Collected:

Date Received:

Field Prep:

Lab ID: L2247717-01 D

Client ID: MW-3

Sample Location: HONEOYE FALLS, NY

Sample Depth:

Matrix: Water
Analytical Method: 1,8260C
Analytical Date: 09/12/22 14:38

Analyst: MV

		Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westboroug	h Lab					
Methylene chloride	ND		ug/l	5.0	1.4	2
1,1-Dichloroethane	ND		ug/l	5.0	1.4	2
Chloroform	ND		ug/l	5.0	1.4	2
Carbon tetrachloride	ND		ug/l	1.0	0.27	2
1,2-Dichloropropane	ND		ug/l	2.0	0.27	2
Dibromochloromethane	ND		ug/l	1.0	0.30	2
1,1,2-Trichloroethane	ND		ug/l	3.0	1.0	2
Tetrachloroethene	5.3		ug/l	1.0	0.36	2
Chlorobenzene	ND		ug/l	5.0	1.4	2
Trichlorofluoromethane	ND		ug/l	5.0	1.4	2
1,2-Dichloroethane	ND		ug/l	1.0	0.26	2
1,1,1-Trichloroethane	8.1		ug/l	5.0	1.4	2
Bromodichloromethane	ND		ug/l	1.0	0.38	2
trans-1,3-Dichloropropene	ND		ug/l	1.0	0.33	2
cis-1,3-Dichloropropene	ND		ug/l	1.0	0.29	2
Bromoform	ND		ug/l	4.0	1.3	2
1,1,2,2-Tetrachloroethane	ND		ug/l	1.0	0.33	2
Benzene	ND		ug/l	1.0	0.32	2
Toluene	ND		ug/l	5.0	1.4	2
Ethylbenzene	ND		ug/l	5.0	1.4	2
Chloromethane	ND		ug/l	5.0	1.4	2
Bromomethane	ND		ug/l	5.0	1.4	2
Vinyl chloride	ND		ug/l	2.0	0.14	2
Chloroethane	ND		ug/l	5.0	1.4	2
1,1-Dichloroethene	3.4		ug/l	1.0	0.34	2
trans-1,2-Dichloroethene	ND		ug/l	5.0	1.4	2
Trichloroethene	290		ug/l	1.0	0.35	2
1,2-Dichlorobenzene	ND		ug/l	5.0	1.4	2



Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

SAMPLE RESULTS

Lab ID: L2247717-01 D Date Collected: 09/01/22 11:18

Client ID: MW-3 Date Received: 09/01/22 Sample Location: HONEOYE FALLS, NY Field Prep: Not Specified

Sample Depth:

No	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1,4-Dichloroberzene ND ug/l 5.0 1.4 2	Volatile Organics by GC/MS - We	estborough Lab					
1.4-Dichlorobenzene ND ug/l 5.0 1.4 2 Methyl tert butyl ether ND ug/l 5.0 1.4 2 p/m-Xylene ND ug/l 5.0 1.4 2 o-Xylene ND ug/l 5.0 1.4 2 o-Xylene ND ug/l 5.0 1.4 2 styrene ND ug/l 5.0 1.4 2 Styrene ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.9 2 Carbon disulfide ND ug/l 10 2.9 2 2-Butanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 5.0 1.4 2 Bromoblytoenzene ND ug/l 5.0 1.4 2 Bromoblytoenzene ND ug/l </td <td>1,3-Dichlorobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>5.0</td> <td>1.4</td> <td>2</td>	1,3-Dichlorobenzene	ND		ug/l	5.0	1.4	2
Methyl tent buyl ether ND ug/l 5.0 1.4 2 p/m-Xylene ND ug/l 5.0 1.4 2 o-Xylene ND ug/l 5.0 1.4 2 cis-1,2-Dichloroethene 24 ug/l 5.0 1.4 2 Styrene ND ug/l 5.0 1.4 2 Styrene ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.0 2 2-Butanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 1-Evanore ND ug/l 5.0 1.4 2 1-2-Dibromoethane ND ug/l 5.0 1.4 2 1-2-Dibromoethane ND ug/l 5.0 1.4 2 1-2-Dibromoethane ND	1,4-Dichlorobenzene	ND			5.0	1.4	2
o-Xylene ND ug/l 5.0 1.4 2 cis-1,2-Dichloroethene 24 ug/l 5.0 1.4 2 Styrene ND ug/l 5.0 1.4 2 Dichlorodifluoromethane ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.9 2 Carbon disulfide ND ug/l 10 2.9 2 2-Butanone ND ug/l 10 3.9 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 5.0 1.4 2 2-Hexanone ND ug/l 5.0 1.4 2 1,2-Distomoethane<	Methyl tert butyl ether	ND			5.0	1.4	2
cis-1,2-Dichloroethene 24 ug/l 5.0 1.4 2 Styrene ND ug/l 5.0 1.4 2 Dichlorodiffuoromethane ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.9 2 Carbon disulfide ND ug/l 10 2.9 2 2-Butanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 5.0 1.4 2 2-Hexanone ND ug/l 5.0 1.4 2 1-Evanone ND ug/l	p/m-Xylene	ND		ug/l	5.0	1.4	2
Styrene ND ug/l 5.0 1.4 2 Dichlorodiffluoromethane ND ug/l 10 2.0 2 Acetone ND ug/l 10 2.9 2 Carbon disulfide ND ug/l 10 2.0 2 2-Butanone ND ug/l 10 3.9 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 5.0 1.4 2 1-2-Dibromo-dance ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND </td <td>o-Xylene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>5.0</td> <td>1.4</td> <td>2</td>	o-Xylene	ND		ug/l	5.0	1.4	2
Dichlorodiffluoromethane	cis-1,2-Dichloroethene	24		ug/l	5.0	1.4	2
Acetone ND ug/l 10 2.9 2 Carbon disulfide ND ug/l 10 2.0 2 2-Butanone ND ug/l 10 2.0 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 Bromochloromethane ND ug/l 5.0 1.4 2 1,2-Dibromo-diane ND ug/l 5.0 1.4 2 n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 P-Isopropytoluene ND ug/l 5.0 1.4 2 1	Styrene	ND			5.0	1.4	2
Carbon disulfide ND ug/l 10 2.0 2 2-Butanone ND ug/l 10 3.9 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 Bromochloromethane ND ug/l 5.0 1.4 2 1,2-Dibromoethane ND ug/l 4.0 1.3 2 n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2	Dichlorodifluoromethane	ND		ug/l	10	2.0	2
2-Butanone ND ug/l 10 3.9 2 4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 Bromochloromethane ND ug/l 5.0 1.4 2 1,2-Dibromoethane ND ug/l 5.0 1.4 2 n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 Isopropyltoluene ND ug/l 5.0 1.4 2 p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2	Acetone	ND		ug/l	10	2.9	2
4-Methyl-2-pentanone ND ug/l 10 2.0 2 2-Hexanone ND ug/l 10 2.0 2 Bromochloromethane ND ug/l 5.0 1.4 2 1,2-Dibromoethane ND ug/l 4.0 1.3 2 n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 sporopylbenzene ND ug/l 5.0 1.4 2 sporopylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2	Carbon disulfide	ND		ug/l	10	2.0	2
2-Hexanone ND ug/l 10 2.0 2	2-Butanone	ND		ug/l	10	3.9	2
Bromochloromethane ND ug/l 5.0 1.4 2	4-Methyl-2-pentanone	ND		ug/l	10	2.0	2
1,2-Dibromoethane ND ug/l 4.0 1.3 2 n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Wethyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 500 120 2 <td>2-Hexanone</td> <td>ND</td> <td></td> <td>ug/l</td> <td>10</td> <td>2.0</td> <td>2</td>	2-Hexanone	ND		ug/l	10	2.0	2
n-Butylbenzene ND ug/l 5.0 1.4 2 sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropylbenzene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 1,4-Dioxane ND ug/l 5.0 1.4 2	Bromochloromethane	ND		ug/l	5.0	1.4	2
sec-Butylbenzene ND ug/l 5.0 1.4 2 1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 5.0 1.4 2 Cyclohexane ND ug/l 4.0 0.47 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2 <td>1,2-Dibromoethane</td> <td>ND</td> <td></td> <td>ug/l</td> <td>4.0</td> <td>1.3</td> <td>2</td>	1,2-Dibromoethane	ND		ug/l	4.0	1.3	2
1,2-Dibromo-3-chloropropane ND ug/l 5.0 1.4 2 Isopropylbenzene ND ug/l 5.0 1.4 2 p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 5.0 1.4 2 Freon-113 ND ug/l 5.0 1.4 2	n-Butylbenzene	ND		ug/l	5.0	1.4	2
Sopropylbenzene ND ug/l 5.0 1.4 2	sec-Butylbenzene	ND		ug/l	5.0	1.4	2
p-Isopropyltoluene ND ug/l 5.0 1.4 2 n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 1,4-Dioxane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	1,2-Dibromo-3-chloropropane	ND		ug/l	5.0	1.4	2
n-Propylbenzene ND ug/l 5.0 1.4 2 1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	Isopropylbenzene	ND		ug/l	5.0	1.4	2
1,2,3-Trichlorobenzene ND ug/l 5.0 1.4 2 1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	p-Isopropyltoluene	ND		ug/l	5.0	1.4	2
1,2,4-Trichlorobenzene ND ug/l 5.0 1.4 2 1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	n-Propylbenzene	ND		ug/l	5.0	1.4	2
1,3,5-Trimethylbenzene ND ug/l 5.0 1.4 2 1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	1,2,3-Trichlorobenzene	ND		ug/l	5.0	1.4	2
1,2,4-Trimethylbenzene ND ug/l 5.0 1.4 2 Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	1,2,4-Trichlorobenzene	ND		ug/l	5.0	1.4	2
Methyl Acetate ND ug/l 4.0 0.47 2 Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	1,3,5-Trimethylbenzene	ND		ug/l	5.0	1.4	2
Cyclohexane ND ug/l 20 0.54 2 1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	1,2,4-Trimethylbenzene	ND		ug/l	5.0	1.4	2
1,4-Dioxane ND ug/l 500 120 2 Freon-113 ND ug/l 5.0 1.4 2	Methyl Acetate	ND		ug/l	4.0	0.47	2
Freon-113 ND ug/l 5.0 1.4 2	Cyclohexane	ND		ug/l	20	0.54	2
-	1,4-Dioxane	ND		ug/l	500	120	2
Methyl cyclohexane ND ug/l 20 0.79 2	Freon-113	ND		ug/l	5.0	1.4	2
	Methyl cyclohexane	ND		ug/l	20	0.79	2

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	100	70-130	
Toluene-d8	100	70-130	
4-Bromofluorobenzene	94	70-130	
Dibromofluoromethane	105	70-130	



09/01/22 11:06

Not Specified

09/01/22

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Lab Number: L2247717

Report Date: 09/16/22

Date Collected:

Date Received:

Field Prep:

SAMPLE RESULTS

Lab ID: L2247717-02 D

Client ID: MW-5

Sample Location: HONEOYE FALLS, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/12/22 15:02

Analyst: MV

1,1-Dichloroethane	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
1,1-Dichloroethane	Volatile Organics by GC/MS - Westbook	rough Lab					
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 Tchloroethane ND ug/l 3.8 1.2 2.5 Chlorobenzene ND ug/l 6.2 1.8 2.5 Trichlorofluoromethane ND ug/l 6.2 1.8 2.5 Trichloroethane ND ug/l 1.2 0.33 2.5 Tchloroptropena ND ug/l 1.2 0.48 2.5 Bromodichloromethane ND ug/l 1.2 0.41 2.5 Bromodichloromethane ND ug/l 1.2 0.41 2.5 Bromodichloromethane ND ug/l 1.2 0.4	Methylene chloride	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 Tetrachloroethane 5.3 ug/l 6.2 1.8 2.5 Chloroberane ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Trichlorofthoroethane ND ug/l 1.2 0.43 2.5 Bromodichloromethane ND ug/l 1.2 0.43 2.5 Bromodichloromethane ND ug/l 1.2 0.48 2.5 Bromodichloromethane ND ug/l 1.2 0.41 2.5 Bromodichloropropene ND ug/l 1.2 0.41 2.5 Bromoform ND ug/l 5.0 <t< td=""><td>1,1-Dichloroethane</td><td>3.4</td><td>J</td><td>ug/l</td><td>6.2</td><td>1.8</td><td>2.5</td></t<>	1,1-Dichloroethane	3.4	J	ug/l	6.2	1.8	2.5
1,2-Dichloropropane ND ug/l 2.5 0.34 2.5 0.5 1,1,2-Trichloroethane ND ug/l 1.2 0.37 2.5 1,1,2-Trichloroethane ND ug/l 3.8 1.2 2.5 1,1,2-Trichloroethane 5.3 ug/l 1.2 0.45 2.5 1,1,2-Trichloroethane ND ug/l 6.2 1.8 2.5 1,2-Dichloroethane ND ug/l 1.2 0.48 2.5 1,2-Dichloropropene ND ug/l 1.2 0.48 2.5 1,3-Dichloropropene ND ug/l 1.2 0.41 2.5 1,3-Dichloropropene ND ug/l 1.2 0.42 2.5 1,1,2,2-Tetrachloroethane ND ug/l 1.2 0.42 2.5 1,1,2,2-Tetrachloroethane ND ug/l 1.2 0.40 2.5 1,1,2,2-Tetrachloroethane ND ug/l 6.2 1.8 2.5 1,1,2,2,2,3,3,3 1.2 1.3 1.3 1,2,2,3,3,3 1.3 1.3 1.3	Chloroform	ND		ug/l	6.2	1.8	2.5
ND	Carbon tetrachloride	ND		ug/l	1.2	0.34	2.5
1,1,2-Trichloroethane	1,2-Dichloropropane	ND		ug/l	2.5	0.34	2.5
Tetrachloroethene 5.3 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 4.0 J ug/l 6.2 1.8 2.5 Bromodichloromethane ND ug/l 1.2 0.48 2.5 Bromodichloropropene ND ug/l 1.2 0.41 2.5 trans-1,3-Dichloropropene ND ug/l 1.2 0.41 2.5 Bromoform ND ug/l 5.0 1.6 2.5 Bromoform ND ug/l 5.0 1.6 2.5 Benzene ND ug/l 6.2 1.8 2.5 Toluene ND ug/l 6.2 1.8 2.5 Ethylbenzene ND ug/l 6.2 1.8	Dibromochloromethane	ND		ug/l	1.2	0.37	2.5
ND	1,1,2-Trichloroethane	ND		ug/l	3.8	1.2	2.5
Trichlorofluoromethane	Tetrachloroethene	5.3		ug/l	1.2	0.45	2.5
1,2-Dichloroethane	Chlorobenzene	ND		ug/l	6.2	1.8	2.5
1,1,1-Trichloroethane	Trichlorofluoromethane	ND		ug/l	6.2	1.8	2.5
Bromodichloromethane ND	1,2-Dichloroethane	ND		ug/l	1.2	0.33	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 Bromoform 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 ND ug/l 2.5 0.18 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 1.2 0.44 2.5 <	1,1,1-Trichloroethane	4.0	J	ug/l	6.2	1.8	2.5
ND	Bromodichloromethane	ND		ug/l	1.2	0.48	2.5
ND ug/l 5.0 1.6 2.5	trans-1,3-Dichloropropene	ND		ug/l	1.2	0.41	2.5
1,1,2,2-Tetrachloroethane	cis-1,3-Dichloropropene	ND		ug/l	1.2	0.36	2.5
ND	Bromoform	ND		ug/l	5.0	1.6	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 ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Trichloroethene DND ug/l 6.2 1.8 2.5 Trichloroethene DND ug/l 6.2 1.8 2.5 Trichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene SOO ug/l 1.2 0.44 2.5	1,1,2,2-Tetrachloroethane	ND		ug/l	1.2	0.42	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 ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Benzene	ND		ug/l	1.2	0.40	2.5
Chloromethane ND ug/l 6.2 1.8 2.5 Bromomethane ND ug/l 6.2 1.8 2.5 Vinyl chloride ND ug/l 2.5 0.18 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Toluene	ND		ug/l	6.2	1.8	2.5
ND ug/l 6.2 1.8 2.5	Ethylbenzene	ND		ug/l	6.2	1.8	2.5
Vinyl chloride ND ug/l 2.5 0.18 2.5 Chloroethane ND ug/l 6.2 1.8 2.5 1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Chloromethane	ND		ug/l	6.2	1.8	2.5
Chloroethane ND ug/l 6.2 1.8 2.5 1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Bromomethane	ND		ug/l	6.2	1.8	2.5
1,1-Dichloroethene 2.4 ug/l 1.2 0.42 2.5 trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Vinyl chloride	ND		ug/l	2.5	0.18	2.5
trans-1,2-Dichloroethene ND ug/l 6.2 1.8 2.5 Trichloroethene 500 ug/l 1.2 0.44 2.5	Chloroethane	ND		ug/l	6.2	1.8	2.5
Trichloroethene 500 ug/l 1.2 0.44 2.5	1,1-Dichloroethene	2.4		ug/l	1.2	0.42	2.5
Ţ	trans-1,2-Dichloroethene	ND		ug/l	6.2	1.8	2.5
1,2-Dichlorobenzene ND ug/l 6.2 1.8 2.5	Trichloroethene	500		ug/l	1.2	0.44	2.5
•	1,2-Dichlorobenzene	ND		ug/l	6.2	1.8	2.5



Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

SAMPLE RESULTS

Lab ID: L2247717-02 D Date Collected: 09/01/22 11:06

Client ID: MW-5 Date Received: 09/01/22 Sample Location: HONEOYE FALLS, NY Field Prep: Not Specified

Sample Depth:

Volatile Organics by GC/MS - Westborough La 1,3-Dichlorobenzene 1,4-Dichlorobenzene Methyl tert butyl ether p/m-Xylene o-Xylene cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone 4-Methyl-2-pentanone	ND ND ND	ug/l	6.2	1.8	
1,4-Dichlorobenzene Methyl tert butyl ether p/m-Xylene o-Xylene cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone	ND		6.2	1.8	
Methyl tert butyl ether p/m-Xylene o-Xylene cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone				1.0	2.5
p/m-Xylene o-Xylene cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone	ND	ug/l	6.2	1.8	2.5
o-Xylene cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone		ug/l	6.2	1.8	2.5
cis-1,2-Dichloroethene Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone	ND	ug/l	6.2	1.8	2.5
Styrene Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone	ND	ug/l	6.2	1.8	2.5
Dichlorodifluoromethane Acetone Carbon disulfide 2-Butanone	140	ug/l	6.2	1.8	2.5
Acetone Carbon disulfide 2-Butanone	ND	ug/l	6.2	1.8	2.5
Carbon disulfide 2-Butanone	ND	ug/l	12	2.5	2.5
2-Butanone	ND	ug/l	12	3.6	2.5
	ND	ug/l	12	2.5	2.5
4-Methyl-2-pentanone	ND	ug/l	12	4.8	2.5
	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
n-Butylbenzene	ND	ug/l	6.2	1.8	2.5
sec-Butylbenzene	ND	ug/l	6.2	1.8	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
p-Isopropyltoluene	ND	ug/l	6.2	1.8	2.5
n-Propylbenzene	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
1,3,5-Trimethylbenzene	ND	ug/l	6.2	1.8	2.5
1,2,4-Trimethylbenzene	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	ua/l	6.2	1.8	2.5
Methyl cyclohexane		ug/l	٥.٢		

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	100	70-130	
Toluene-d8	102	70-130	
4-Bromofluorobenzene	95	70-130	
Dibromofluoromethane	110	70-130	



09/01/22 11:43

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

SAMPLE RESULTS

Lab Number: L2247717

Report Date: 09/16/22

Lab ID: L2247717-03 D Date Collected:

Client ID: Date Received: 09/01/22 MW-201D Field Prep: Sample Location: HONEOYE FALLS, NY Not Specified

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/12/22 15:25

Analyst: MV

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
Volatile Organics by GC/MS - Wes	stborough Lab						
Methylene chloride	ND		ug/l	120	35.	50	
1,1-Dichloroethane	ND		ug/l	120	35.	50	
Chloroform	ND		ug/l	120	35.	50	
Carbon tetrachloride	ND		ug/l	25	6.7	50	
1,2-Dichloropropane	ND		ug/l	50	6.8	50	
Dibromochloromethane	ND		ug/l	25	7.4	50	
1,1,2-Trichloroethane	ND		ug/l	75	25.	50	
Tetrachloroethene	54		ug/l	25	9.0	50	
Chlorobenzene	ND		ug/l	120	35.	50	
Trichlorofluoromethane	ND		ug/l	120	35.	50	
1,2-Dichloroethane	ND		ug/l	25	6.6	50	
1,1,1-Trichloroethane	160		ug/l	120	35.	50	
Bromodichloromethane	ND		ug/l	25	9.6	50	
trans-1,3-Dichloropropene	ND		ug/l	25	8.2	50	
cis-1,3-Dichloropropene	ND		ug/l	25	7.2	50	
Bromoform	ND		ug/l	100	32.	50	
1,1,2,2-Tetrachloroethane	ND		ug/l	25	8.4	50	
Benzene	ND		ug/l	25	8.0	50	
Toluene	ND		ug/l	120	35.	50	
Ethylbenzene	ND		ug/l	120	35.	50	
Chloromethane	ND		ug/l	120	35.	50	
Bromomethane	ND		ug/l	120	35.	50	
Vinyl chloride	ND		ug/l	50	3.6	50	
Chloroethane	ND		ug/l	120	35.	50	
1,1-Dichloroethene	70		ug/l	25	8.4	50	
trans-1,2-Dichloroethene	ND		ug/l	120	35.	50	
Trichloroethene	6000		ug/l	25	8.8	50	
1,2-Dichlorobenzene	ND		ug/l	120	35.	50	

MDL

Dilution Factor

Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

SAMPLE RESULTS

Lab ID: L2247717-03 D Date Collected: 09/01/22 11:43

Client ID: MW-201D Date Received: 09/01/22 Sample Location: HONEOYE FALLS, NY Field Prep: Not Specified

Qualifier

Units

RL

Result

Sample Depth:

Parameter

- urameter						
Volatile Organics by GC/MS - Wes	stborough Lab					
1,3-Dichlorobenzene	ND	ug/l	120	35.	50	
1,4-Dichlorobenzene	ND	ug/l	120	35.	50	
Methyl tert butyl ether	ND	ug/l	120	35.	50	
p/m-Xylene	ND	ug/l	120	35.	50	
o-Xylene	ND	ug/l	120	35.	50	
cis-1,2-Dichloroethene	1200	ug/l	120	35.	50	
Styrene	ND	ug/l	120	35.	50	
Dichlorodifluoromethane	ND	ug/l	250	50.	50	
Acetone	ND	ug/l	250	73.	50	
Carbon disulfide	ND	ug/l	250	50.	50	
2-Butanone	ND	ug/l	250	97.	50	
4-Methyl-2-pentanone	ND	ug/l	250	50.	50	
2-Hexanone	ND	ug/l	250	50.	50	
Bromochloromethane	ND	ug/l	120	35.	50	
1,2-Dibromoethane	ND	ug/l	100	32.	50	
n-Butylbenzene	ND	ug/l	120	35.	50	
sec-Butylbenzene	ND	ug/l	120	35.	50	
1,2-Dibromo-3-chloropropane	ND	ug/l	120	35.	50	
Isopropylbenzene	ND	ug/l	120	35.	50	
p-Isopropyltoluene	ND	ug/l	120	35.	50	
n-Propylbenzene	ND	ug/l	120	35.	50	
1,2,3-Trichlorobenzene	ND	ug/l	120	35.	50	
1,2,4-Trichlorobenzene	ND	ug/l	120	35.	50	
1,3,5-Trimethylbenzene	ND	ug/l	120	35.	50	
1,2,4-Trimethylbenzene	ND	ug/l	120	35.	50	
Methyl Acetate	ND	ug/l	100	12.	50	
Cyclohexane	ND	ug/l	500	14.	50	
1,4-Dioxane	ND	ug/l	12000	3000	50	
Freon-113	ND	ug/l	120	35.	50	
Methyl cyclohexane	ND	ug/l	500	20.	50	

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	103	70-130	
Toluene-d8	102	70-130	
4-Bromofluorobenzene	96	70-130	
Dibromofluoromethane	110	70-130	



Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

SAMPLE RESULTS

Lab Number: L2247717

Report Date: 09/16/22

Lab ID: L2247717-04

Client ID: SUPPLY WELL

Sample Location: HONEOYE FALLS, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/12/22 16:12

Analyst: MV

Date Collected:	09/01/22 11:22
Date Received:	09/01/22
Field Prep:	Not Specified

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 Chrorotetrachioride ND ug/l 0.50 0.13 1 L2-Dichloropropane ND ug/l 0.50 0.13 1 L2-Dichloropropane ND ug/l 0.50 0.15 1 Dibromochloromethane ND ug/l 0.50 0.15 1 1,1,2-Trichloroethane ND ug/l 0.50 0.18 1 1,2-Dichloroethane ND ug/l 0.50 0.18 1 1,2-Dichloroethane ND ug/l 0.50 0.13 1 1,1-1-Trichloroethane ND ug/l 0.50 0.13 1 1,1-1-Trichloroethane ND ug/l 0.50 0.19 1 1,1-1-Trichloroethane	Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor	
1,1-Dichloroethane ND ug/l 2.5 0.70 1	Volatile Organics by GC/MS - Wes	stborough Lab						
Chloroform ND ugfl 2.5 0.70 1 Carbon tetrachloride ND ugfl 0.50 0.13 1 1,2-Dichloropropane ND ugfl 1.0 0.14 1 Dibromochloromethane ND ugfl 0.50 0.15 1 1,1,2-Trichloroethane ND ugfl 0.50 0.18 1 1,1,2-Trichloroethane ND ugfl 0.50 0.18 1 Chlorobenzene ND ugfl 2.5 0.70 1 Trichlorofluoromethane ND ugfl 2.5 0.70 1 1,2-Dichloropethane ND ugfl 2.5 0.70 1 1,1-1-Trichloroethane ND ugfl 0.50 0.13 1 Bromodichloromethane ND ugfl 0.50 0.19 1 trans-1,3-Dichloropropene ND ugfl 0.50 0.16 1 Bromodichloromethane ND ugfl 0.50	Methylene chloride	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 Tetrachloroethane 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 0.50 0.13 1 Bromodichloromethane ND ug/l 0.50 0.19 1 Bromodichloropropene ND ug/l 0.50 0.16 1 cis-1,3-Dichloropropene ND ug/l 0.50 0.14 1 Bromodorm ND ug/l 0.50 <t< td=""><td>1,1-Dichloroethane</td><td>ND</td><td></td><td>ug/l</td><td>2.5</td><td>0.70</td><td>1</td><td></td></t<>	1,1-Dichloroethane	ND		ug/l	2.5	0.70	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 Tetrachloroethane 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 0.50 0.13 1 Bromodichloromethane ND ug/l 0.50 0.13 1 Bromodichloropropene ND ug/l 0.50 0.16 1 cis-1,3-Dichloropropene ND ug/l 0.50 0.16 1 sis-1,3-Dichloropropene ND ug/l 0.50 0.16 1 Benzene ND ug/l 0.50 <	Chloroform	ND		ug/l	2.5	0.70	1	
Dibromochloromethane ND ug/l 0.50 0.15 1 1,1,2-Trichloroethane ND ug/l 1.5 0.50 1 Tetrachloroethane ND ug/l 0.50 0.18 1 Chlorobenzene ND ug/l 2.5 0.70 1 Trichloroffluoromethane 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 0.50 0.13 1 Bromodichloromethane ND ug/l 0.50 0.13 1 Bromodichloromethane ND ug/l 0.50 0.16 1 trans-1,3-Dichloropropene ND ug/l 0.50 0.16 1 Bromoform ND ug/l 0.50 0.16 1 Toluene ND ug/l 0.50 0.17 1 Benzene ND ug/l 2.5 0.70 <t< td=""><td>Carbon tetrachloride</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>0.13</td><td>1</td><td></td></t<>	Carbon tetrachloride	ND		ug/l	0.50	0.13	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 0.13 1 Bromodichloromethane ND ug/l 0.50 0.19 1 trans-1,3-Dichloropropene ND ug/l 0.50 0.16 1 ticis-1,3-Dichloropropene ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.16 1 1,1,2,2-Tetrachloroethane ND ug/l 2.5 <	1,2-Dichloropropane	ND		ug/l	1.0	0.14	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 0.50 0.19 1 Bromodichloromethane ND ug/l 0.50 0.19 1 It trans-1,3-Dichloropropene ND ug/l 0.50 0.16 1 cis-1,3-Dichloropropene ND ug/l 0.50 0.16 1 strans-1,3-Dichloropropene ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.14 1 Bromoform 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 <t< td=""><td>Dibromochloromethane</td><td>ND</td><td></td><td>ug/l</td><td>0.50</td><td>0.15</td><td>1</td><td></td></t<>	Dibromochloromethane	ND		ug/l	0.50	0.15	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 0.50 0.14 1 Bromoform ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.17 1 Benzene ND ug/l 0.50 0.16 1 1,1,2,2-Tetrachloroethane ND ug/l 2.5 0.70 1 Ethylbenzene ND ug/l 2.5 0.70 1	1,1,2-Trichloroethane	ND		ug/l	1.5	0.50	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 dis-1,3-Dichloropropene ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.14 1 Bromoform ND ug/l 0.50 0.14 1 1,1,2,2-Tetrachloroethane ND ug/l 0.50 0.16 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	Tetrachloroethene	ND		ug/l	0.50	0.18	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 sis-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 Stromomethane ND ug/l 2.5 0.70 1 Vinyl chloride ND ug/l 2.5 0.70 1 <td>Chlorobenzene</td> <td>ND</td> <td></td> <td>ug/l</td> <td>2.5</td> <td>0.70</td> <td>1</td> <td></td>	Chlorobenzene	ND		ug/l	2.5	0.70	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 2.5 0.70 1 Chloroethane ND ug/l 2.5 0.70 1 <t< td=""><td>Trichlorofluoromethane</td><td>ND</td><td></td><td>ug/l</td><td>2.5</td><td>0.70</td><td>1</td><td></td></t<>	Trichlorofluoromethane	ND		ug/l	2.5	0.70	1	
Bromodichloromethane ND	1,2-Dichloroethane	ND		ug/l	0.50	0.13	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 2.5 0.70 1 Chloroethane ND ug/l 2.5 0.70 1 1,1-Dichloroethene 0.27 J ug/l 2.5 0.70 1 1,1-Dichloroethene ND ug/l 2.5 0.70 1	1,1,1-Trichloroethane	ND		ug/l	2.5	0.70	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Bromodichloromethane	ND		ug/l	0.50	0.19	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 2.5 0.70 1 Chloroethane ND ug/l 2.5 0.70 1 1,1-Dichloroethene 0.27 J ug/l 2.5 0.70 1 1,1-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	trans-1,3-Dichloropropene	ND		ug/l	0.50	0.16	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	cis-1,3-Dichloropropene	ND		ug/l	0.50	0.14	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Bromoform	ND		ug/l	2.0	0.65	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	1,1,2,2-Tetrachloroethane	ND		ug/l	0.50	0.17	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Benzene	ND		ug/l	0.50	0.16	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Toluene	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Ethylbenzene	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 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Chloromethane	ND		ug/l	2.5	0.70	1	
Chloroethane ND ug/l 2.5 0.70 1 1,1-Dichloroethene 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Bromomethane	ND		ug/l	2.5	0.70	1	
1,1-Dichloroethene 0.27 J ug/l 0.50 0.17 1 trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Vinyl chloride	ND		ug/l	1.0	0.07	1	
trans-1,2-Dichloroethene ND ug/l 2.5 0.70 1 Trichloroethene 12 ug/l 0.50 0.18 1	Chloroethane	ND		ug/l	2.5	0.70	1	
Trichloroethene 12 ug/l 0.50 0.18 1	1,1-Dichloroethene	0.27	J	ug/l	0.50	0.17	1	
	trans-1,2-Dichloroethene	ND		ug/l	2.5	0.70	1	
A O Dichlorate areas	Trichloroethene	12		ug/l	0.50	0.18	1	
1,z-Dichlorobenzene ND ug/l 2.5 0.70 1	1,2-Dichlorobenzene	ND		ug/l	2.5	0.70	1	



Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

SAMPLE RESULTS

Lab ID: L2247717-04 Date Collected: 09/01/22 11:22

Client ID: SUPPLY WELL Date Received: 09/01/22
Sample Location: HONEOYE FALLS, NY Field Prep: Not Specified

Sample Depth:

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westl	borough Lab					
1,3-Dichlorobenzene	ND		ug/l	2.5	0.70	1
1,4-Dichlorobenzene	ND		ug/l	2.5	0.70	1
Methyl tert butyl ether	ND		ug/l	2.5	0.70	1
p/m-Xylene	ND		ug/l	2.5	0.70	1
o-Xylene	ND		ug/l	2.5	0.70	1
cis-1,2-Dichloroethene	1.1	J	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
n-Butylbenzene	ND		ug/l	2.5	0.70	1
sec-Butylbenzene	ND		ug/l	2.5	0.70	1
1,2-Dibromo-3-chloropropane	ND		ug/l	2.5	0.70	1
Isopropylbenzene	ND		ug/l	2.5	0.70	1
p-Isopropyltoluene	ND		ug/l	2.5	0.70	1
n-Propylbenzene	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
1,3,5-Trimethylbenzene	ND		ug/l	2.5	0.70	1
1,2,4-Trimethylbenzene	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	105	70-130	
Toluene-d8	101	70-130	
4-Bromofluorobenzene	94	70-130	
Dibromofluoromethane	108	70-130	



09/01/22 00:00

Not Specified

09/01/22

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

SAMPLE RESULTS

Lab Number: L2247717

Report Date: 09/16/22

Date Collected:

Date Received:

Field Prep:

Lab ID: L2247717-05

Client ID: TRIP BLANK

Sample Location: HONEOYE FALLS, NY

Sample Depth:

Matrix: Water Analytical Method: 1,8260C Analytical Date: 09/12/22 09:13

Analyst: MKS

Parameter	Result	Qualifier	Units	RL	MDL	Dilution Factor
Volatile Organics by GC/MS - Westbook	ough 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: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

SAMPLE RESULTS

Lab ID: L2247717-05 Date Collected: 09/01/22 00:00

Client ID: TRIP BLANK Date Received: 09/01/22 Sample Location: HONEOYE FALLS, NY Field Prep: Not Specified

Sample Depth:

1,3-Dichlorobenzene	2.5 2.5 2.5 2.5 2.5 2.5 2.5 2.5 5.0	0.70 0.70 0.70 0.70 0.70 0.70	1 1 1 1
1,4-Dichlorobenzene ND ug/l Methyl tert butyl ether ND ug/l p/m-Xylene ND ug/l c-Xylene ND ug/l cis-1,2-Dichloroethene ND ug/l Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l lsopropyltenzene ND ug/l lsopropyltouene ND ug/l n-Propyltenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5 2.5 2.5 2.5 2.5 2.5 2.5 5.0	0.70 0.70 0.70 0.70 0.70	1 1 1
1,4-Dichlorobenzene ND ug/l Methyl tert butyl ether ND ug/l p/m-Xylene ND ug/l o-Xylene ND ug/l cis-1,2-Dichloroethene ND ug/l Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropylbenzene ND ug/l n-Propylbenzene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene	2.5 2.5 2.5 2.5 2.5 2.5 5.0	0.70 0.70 0.70 0.70	1 1 1
Methyl tert butyl ether ND ug/l p/m-Xylene ND ug/l o-Xylene ND ug/l cis-1,2-Dichloroethene ND ug/l Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropyltenzene ND ug/l n-Propylbenzene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5 2.5 2.5 2.5 2.5 5.0	0.70 0.70 0.70	1
p/m-Xylene ND ug/l o-Xylene ND ug/l cis-1,2-Dichloroethene ND ug/l Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5 2.5 2.5 5.0	0.70 0.70	1
cis-1,2-Dichloroethene ND ug/l Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l 1sopropyltenzene ND ug/l n-Propylbenzene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5 2.5 5.0	0.70	
Styrene ND ug/l Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5 5.0		
Dichlorodifluoromethane ND ug/l Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	0.70	1
Acetone ND ug/l Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l p-Isopropylbenzene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l		0.70	1
Carbon disulfide ND ug/l 2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	1.0	1
2-Butanone ND ug/l 4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l		1.5	1
4-Methyl-2-pentanone ND ug/l 2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l lsopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	1.0	1
2-Hexanone ND ug/l Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	1.9	1
Bromochloromethane ND ug/l 1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	1.0	1
1,2-Dibromoethane ND ug/l n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	5.0	1.0	1
n-Butylbenzene ND ug/l sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
sec-Butylbenzene ND ug/l 1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.0	0.65	1
1,2-Dibromo-3-chloropropane ND ug/l Isopropylbenzene ND ug/l p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
Isopropylbenzene	2.5	0.70	1
p-Isopropyltoluene ND ug/l n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
n-Propylbenzene ND ug/l 1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
1,2,3-Trichlorobenzene ND ug/l 1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
1,2,4-Trichlorobenzene ND ug/l	2.5	0.70	1
	2.5	0.70	1
1,3,5-Trimethylbenzene ND ug/l	2.5	0.70	1
	2.5	0.70	1
1,2,4-Trimethylbenzene ND ug/l	2.5	0.70	1
Methyl Acetate ND ug/l	2.0	0.23	1
Cyclohexane ND ug/l		0.27	1
1,4-Dioxane ND ug/l	10	61.	1
Freon-113 ND ug/l	10 250	0.70	1
Methyl cyclohexane ND ug/l			1

Surrogate	% Recovery	Acceptance Qualifier Criteria	
1,2-Dichloroethane-d4	98	70-130	
Toluene-d8	101	70-130	
4-Bromofluorobenzene	98	70-130	
Dibromofluoromethane	105	70-130	



Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/12/22 08:28

Analyst: MV

arameter	Result	Qualifier Units	s RL	MDL
olatile Organics by GC/MS	- Westborough Lab	for sample(s):	01-05 Batch:	WG1686832-5
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: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

Method Blank Analysis Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/12/22 08:28

Analyst: MV

arameter	Result	Qualifier Units	RL	MDL
olatile Organics by GC/MS - \	Westborough Lab	for sample(s):	01-05 Batch:	WG1686832-5
1,4-Dichlorobenzene	ND	ug/l	2.5	0.70
Methyl tert butyl ether	ND	ug/l	2.5	0.70
p/m-Xylene	ND	ug/l	2.5	0.70
o-Xylene	ND	ug/l	2.5	0.70
cis-1,2-Dichloroethene	ND	ug/l	2.5	0.70
Styrene	ND	ug/l	2.5	0.70
Dichlorodifluoromethane	ND	ug/l	5.0	1.0
Acetone	ND	ug/l	5.0	1.5
Carbon disulfide	ND	ug/l	5.0	1.0
2-Butanone	ND	ug/l	5.0	1.9
4-Methyl-2-pentanone	ND	ug/l	5.0	1.0
2-Hexanone	ND	ug/l	5.0	1.0
Bromochloromethane	ND	ug/l	2.5	0.70
1,2-Dibromoethane	ND	ug/l	2.0	0.65
n-Butylbenzene	ND	ug/l	2.5	0.70
sec-Butylbenzene	ND	ug/l	2.5	0.70
1,2-Dibromo-3-chloropropane	ND	ug/l	2.5	0.70
Isopropylbenzene	ND	ug/l	2.5	0.70
p-Isopropyltoluene	ND	ug/l	2.5	0.70
n-Propylbenzene	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
1,3,5-Trimethylbenzene	ND	ug/l	2.5	0.70
1,2,4-Trimethylbenzene	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: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 **Report Date:** 09/16/22

Method Blank Analysis
Batch Quality Control

Analytical Method: 1,8260C Analytical Date: 09/12/22 08:28

Analyst: MV

Parameter Result Qualifier Units RL MDL

Volatile Organics by GC/MS - Westborough Lab for sample(s): 01-05 Batch: WG1686832-5

		Acceptance			
Surrogate	%Recovery	Qualifier C	riteria		
1,2-Dichloroethane-d4	101	70)-130		
Toluene-d8	101	70)-130		
4-Bromofluorobenzene	99	70)-130		
Dibromofluoromethane	108	70)-130		



Lab Control Sample Analysis Batch Quality Control

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Lab Number: L2247717

Report Date: 09/16/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery		%Recovery Limits	RPD	RPD Qual Limits	
/olatile Organics by GC/MS - Westborough	Lab Associated	sample(s):	01-05 Batch:	WG1686832-3	WG1686832-4			
Methylene chloride	97		95		70-130	2	20	
1,1-Dichloroethane	98		96		70-130	2	20	
Chloroform	97		98		70-130	1	20	
Carbon tetrachloride	98		98		63-132	0	20	
1,2-Dichloropropane	96		98		70-130	2	20	
Dibromochloromethane	91		100		63-130	9	20	
1,1,2-Trichloroethane	91		95		70-130	4	20	
Tetrachloroethene	99		97		70-130	2	20	
Chlorobenzene	96		96		75-130	0	20	
Trichlorofluoromethane	120		120		62-150	0	20	
1,2-Dichloroethane	92		96		70-130	4	20	
1,1,1-Trichloroethane	100		99		67-130	1	20	
Bromodichloromethane	95		97		67-130	2	20	
trans-1,3-Dichloropropene	96		96		70-130	0	20	
cis-1,3-Dichloropropene	94		96		70-130	2	20	
Bromoform	80		86		54-136	7	20	
1,1,2,2-Tetrachloroethane	80		88		67-130	10	20	
Benzene	98		97		70-130	1	20	
Toluene	97		96		70-130	1	20	
Ethylbenzene	95		94		70-130	1	20	
Chloromethane	90		87		64-130	3	20	
Bromomethane	100		100		39-139	0	20	
Vinyl chloride	100		97		55-140	3	20	



Lab Control Sample Analysis Batch Quality Control

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Lab Number: L2247717

Report Date: 09/16/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	Qual	%Recovery Limits	RPD	RPD Qual Limits
Volatile Organics by GC/MS - Westborough	Lab Associated	sample(s):	01-05 Batch:	WG1686832-3	WG1686832-4		
Chloroethane	140	Q	140	Q	55-138	0	20
1,1-Dichloroethene	96		94		61-145	2	20
trans-1,2-Dichloroethene	100		97		70-130	3	20
Trichloroethene	93		92		70-130	1	20
1,2-Dichlorobenzene	89		92		70-130	3	20
1,3-Dichlorobenzene	91		95		70-130	4	20
1,4-Dichlorobenzene	90		91		70-130	1	20
Methyl tert butyl ether	81		87		63-130	7	20
p/m-Xylene	95		95		70-130	0	20
o-Xylene	90		90		70-130	0	20
cis-1,2-Dichloroethene	98		98		70-130	0	20
Styrene	95		95		70-130	0	20
Dichlorodifluoromethane	66		63		36-147	5	20
Acetone	76		80		58-148	5	20
Carbon disulfide	97		93		51-130	4	20
2-Butanone	82		85		63-138	4	20
4-Methyl-2-pentanone	80		83		59-130	4	20
2-Hexanone	71		74		57-130	4	20
Bromochloromethane	99		100		70-130	1	20
1,2-Dibromoethane	91		95		70-130	4	20
n-Butylbenzene	93		93		53-136	0	20
sec-Butylbenzene	94		95		70-130	1	20
1,2-Dibromo-3-chloropropane	74		86		41-144	15	20



Lab Control Sample Analysis Batch Quality Control

Project Name: ALCO HONEOYE GWM

Project Number: B0672-022-001

Lab Number: L2247717

Report Date: 09/16/22

Parameter	LCS %Recovery	Qual	LCSD %Recovery	' Qual	%Recovery Limits	RPD	Qual	RPD Limits
Volatile Organics by GC/MS - Westborough L	ab Associated	sample(s):	01-05 Batch:	WG1686832-3	WG1686832-4			
Isopropylbenzene	92		94		70-130	2		20
p-Isopropyltoluene	93		93		70-130	0		20
n-Propylbenzene	93		96		69-130	3		20
1,2,3-Trichlorobenzene	83		92		70-130	10		20
1,2,4-Trichlorobenzene	85		93		70-130	9		20
1,3,5-Trimethylbenzene	92		92		64-130	0		20
1,2,4-Trimethylbenzene	90		91		70-130	1		20
Methyl Acetate	78		82		70-130	5		20
Cyclohexane	95		93		70-130	2		20
1,4-Dioxane	122		128		56-162	5		20
Freon-113	98		97		70-130	1		20
Methyl cyclohexane	92		92		70-130	0		20

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



Serial_No:09162213:57 Lab Number: L2247717

Project Name: ALCO HONEOYE GWM **Project Number:** B0672-022-001

YES

Report Date: 09/16/22

Sample Receipt and Container Information

Were project specific reporting limits specified?

Cooler Information

Custody Seal Cooler

Α Absent

Container Information			Initial	Final	Temp			Frozen		
Container ID	Container Type	Cooler	рН	pН	deg C	Pres	Seal	Date/Time	Analysis(*)	
L2247717-01A	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-01B	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-01C	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-02A	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-02B	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-02C	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-03A	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-03B	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-03C	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-04A	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-04B	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-04C	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-05A	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	
L2247717-05B	Vial HCl preserved	Α	NA		4.1	Υ	Absent		NYTCL-8260-R2(14)	



Project Name: Lab Number: ALCO HONEOYE GWM L2247717 **Project Number:** B0672-022-001 **Report Date:** 09/16/22

GLOSSARY

Acronyms

LCSD

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.

 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

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

MDI - 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.

TEO - 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:ALCO HONEOYE GWMLab Number:L2247717Project Number:B0672-022-001Report Date:09/16/22

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, Benza(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.
- 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 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).
- Co-elution: The target analyte co-elutes with a known lab standard (i.e. surrogate, internal standards, etc.) for co-extracted analyses.
- 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.
- 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

Report Format: DU Report with 'J' Qualifiers



Project Name:ALCO HONEOYE GWMLab Number:L2247717Project Number:B0672-022-001Report Date:09/16/22

Data Qualifiers

Identified Compounds (TICs).

- $\label{eq:main_main_model} \textbf{M} \qquad \text{-Reporting Limit (RL) exceeds the MCP CAM Reporting Limit for this analyte.}$
- 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.
- ${f P}$ 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.
- The surrogate associated with this target analyte has a recovery outside the QC acceptance limits. (Applicable to MassDEP DW Compliance samples only.)
- 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.)

Report Format: DU Report with 'J' Qualifiers



Serial_No:09162213:57

Project Name: ALCO HONEOYE GWM Lab Number: L2247717

Project Number: B0672-022-001 Report Date: 09/16/22

REFERENCES

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.



Serial_No:09162213:57

Alpha Analytical, Inc. Facility: Company-wide

Department: Quality Assurance

Title: Certificate/Approval Program Summary

ID No.:17873 Revision 19

Published Date: 4/2/2021 1:14:23 PM

Page 1 of 1

Certification Information

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

Westborough Facility

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

EPA 625/625.1: alpha-Terpineol

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

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

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

Mansfield Facility

SM 2540D: TSS

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

EPA TO-15: Halothane, 2,4,4-Trimethyl-2-pentene, 2,4,4-Trimethyl-1-pentene, Thiophene, 2-Methylthiophene,

3-Methylthiophene, 2-Ethylthiophene, 1,2,3-Trimethylbenzene, Indan, Indene, 1,2,4,5-Tetramethylbenzene, Benzothiophene, 1-Methylnaphthalene.

Biological Tissue Matrix: EPA 3050B

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

Westborough Facility:

Drinking Water

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

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

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

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

Non-Potable Water

SM4500H,B, EPA 120.1, SM2510B, SM2540C, SM2320B, SM4500CL-E, SM4500F-BC, SM4500NH3-BH: Ammonia-N and Kjeldahl-N, EPA 350.1: Ammonia-N, 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 II, Endosulfan II, Endosulfan sulfate, Endrin, Endrin Aldehyde, Heptachlor, Heptachlor Epoxide, PCBs

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

Microbiology: SM9223B-Colilert-QT; Enterolert-QT, SM9221E, EPA 1600, EPA 1603, SM9222D.

Mansfield Facility:

Drinking Water

EPA 200.7: Al, Ba, Cd, Cr, Cu, Fe, Mn, Ni, Na, Ag, Ca, Zn. EPA 200.8: Al, Sb, As, Ba, Be, Cd, Cr, Cu, Pb, Mn, Ni, Se, Ag, TL, Zn. EPA 245.1 Hg. EPA 522, EPA 537.1.

Non-Potable Water

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

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

EPA 245.1 Hg

SM2340B

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

Document Type: Form

Pre-Qualtrax Document ID: 08-113

Westborough, MA 01581	NEW YORK CHAIN OF CUSTODY Mansfield, MA 02048	Service Centers Mahwah, NJ 07430: 35 Whitne Albany, NY 12205: 14 Walker V Tonawanda, NY 14150: 275 Co	Way ooper Av	e, Suite 1	_		e of /		Date Rec'd in Lab	9/	ALPHA Job # L2247717 Billing Information		
8 Walkup Dr. TEL: 508-898-9220 FAX: 508-898-9193	320 Forbes Blvd TEL 508-822-9300 FAX: 508-822-3288	Project Name: Alce Project Location: How	er o	eye	fells	GWM.			ASP-A EQuIS (1 Fil	_	ASP-B EQuIS (4 File)	Same as Client Info	
Client Information	-	Project# 13067	2-1	022	-001			-	Other	smant		Disposal Site Information	
Address: 2542 Ho	ENLY Turpiles	ALPHAQuote #:	roject #		tann				alory Requir NY TOGS AWQ Standard NY Restricted	fs 🗍	NY Part 375 NY CP-51 Other	Please identify below location of applicable disposal facilities. Disposal Facility:	
Phone: (7/6) & Fax: Email: TRANSMENT		Turn-Around Time Standard Rush (only if pre approved			Due Date				NY Unrestricte NYC Sewer Di	d Use	Other	NJ NY Other:	
	been previously analyz							ANAL	YSIS			Sample Filtration	T
Other project specifi	c requirements/comm	nents:						+CP-51 8260				Preservation Lab to do	1 B - B - 1
ALPHA Lab ID	S	ample ID				Sample			Tat				1
-	(Lab Use Only)		Date Time		Matrix Initials		F					B	
47717 01	mw-3		19/	122	1115	Witer	TAG	X					3
202	MW-5				1106	1		×					3
-03	Mw - 2011				1143			×					3
704	Supply w	ال	-	S	1122			X					3
705	Tric Bla	ile				1		X					2
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Preservative Code: A = None B = HCl C = HNO ₁ D = H ₂ SO ₄ E = NaOH	Container Code P = Plastic A = Amber Glass V = Vial G = Glass B = Bacteria Cup	Westboro: Certification No: MA935 Mansfield: Certification No: MA015				Container Type Preservative		Q I				Please print clearly, legibly and completely. Samples can not be logged in and turnaround time clock will not start until any ambiguities are	
F = MeOH G = NaHSO ₄ H = Na ₂ S ₂ O ₅ K/E = Zn Ac/NaOH O = Other Form No. 01-25 HC (rev. 3	C = Cube O = Other E = Encore D = BOD Bottle	1 17				/Time 15/5 /632	3	Received By			Pate/Time <u>22 </u>	resolved. BY EXECUTING THIS COC, THE CLIENT HAS READ AND AGREES	

HISTORIC GROUNDWATER MONITORING RESULTS





SUMMARY OF HISTORIC ON-SITE GROUNDWATER ANALYTICAL RESULTS

Enarc-O Machine Products, Inc. Lima, New York NYSDEC Registry No. 8-26-011

	DATE	COMPOUND							
WELL		1,1,1-TCA	1,1-DCE	cis-1,2-DCE	TCE	PCE	Toluene	1,1-DCA	Total VOCs
	25-Feb-91								0
	14-Jul-94	130	14 J	30 J	1100	17 J			1291
	2-Nov-94	250		51 J	3200	23 J			3524
	14-Apr-95	190	12	98	2500	22			2822
	23-Aug-95	47	4 J	22	510	10			593
	27-Oct-99	525			8650				9175
	8-Feb-00	365			5250				5615
	27-Apr-00	43.2			585				628
	25-Jul-00	121			1780				1901
	19-Oct-00	502		315	6830				7647
	21-Dec-00	57.8		103	1020				1181
	28-Feb-01			154	1630				1784
	19-Apr-01	167		174	2950				3291
	25-Oct-01	382		746	7210				8338
	11-Apr-02			105	1860				1965
	29-Oct-02	464		347	6390				7201
	29-Apr-03	250		268	4050				4568
MW-3	27-Oct-03	285		288	5720				6293
	29-Apr-04	261		152	3550				3963
	28-Oct-04	390		504	8430				9324
	12-Feb-07	97	18	440	1800				2355
	15-Aug-07	24		45	440	4.7 J			514
	13-Mar-08	38	10	210	930 D	4.5 J			1193
	20-Nov-08	22	5.9	63	490	6			587
	4-Feb-10	ND	ND	140	830	ND	ND		970
	1-May-11	11	ND	40	300	ND	ND		351
	29-Sep-12	ND	ND	24	300	ND	ND		324
	13-Nov-13	7.3	ND	12	180	ND	ND		199
	20-Feb-15	11	ND	95 D	610 D	8.4	ND		724
	31-May-16	ND	ND	49	360	ND	ND		409
	24-Aug-17	13	10	19	260	4.9	ND		307
	20-Nov-18	ND	ND	7.9	120	ND	ND		128
	13-Jan-20	10	7.4	24	380	6.3	ND		428
	24-May-21	9.8	4.2	45	380	6.4	ND		445
	1-Sep-22	8.1	3.4	24	290	5.3	ND	ND	330.8
	7-Jan-91								ND
	25-Feb-91	22.1		50	540				ND
	14-Jul-94 2-Nov-94	23 J 55	5 J	58 72	510 1100	9 J			591 1241
	14-Apr-95	15		63	400	4 J			482
	23-Aug-95	73	7 J	67	540	7 J			694
	27-Oct-99	33	7		657	6			703
MW-5	8-Feb-00	8.5		27.4 J	170				179
	27-Apr-00	5.24			161 1120				166
	25-Jul-00 19-Oct-00	47.8 8.6	2.01	30.1	199				1168 240
	21-Dec-00	7.14	2.01	36.1	163				206
	28-Feb-01	2.03		29.3	78.3				110
	19-Apr-01	2.4	2.46	49.3	114				168
	25-Oct-01	35.6		139	758				933
	11-Apr-02 29-Oct-02	4.8 45		89 158	191 953	10.8			285 1167
	29-Oct-02 29-Apr-03	6.17	2.78	84.8	953 222	10.0			316
	27-Oct-03	28.5	2.70	90.2	698				817
	29-Apr-04	4.01		71.7	178				254
	28-Oct-04	88	24	324	2300				2736
l	12-Feb-07	42	20	490	970		l		1522



SUMMARY OF HISTORIC ON-SITE GROUNDWATER ANALYTICAL RESULTS

Enarc-O Machine Products, Inc. Lima, New York NYSDEC Registry No. 8-26-011

	NYSDEC Registry No. 8-26-011 COMPOUND								
WELL	DATE								Total VOCs
MW-5	15-Aug-07	1,1,1-TCA 28	1,1-DCE 11 J	cis-1,2-DCE 360	1300	PCE	Toluene	1,1-DCA	
ivivv-5	12-Mar-08	1.3	21 J	27	88	0.51 J			1699 138
	20-Nov-08	38	15	390	1400	13			1856
	4-Feb-10	ND	ND	110	290	ND	ND		400
	1-May-11	ND	ND	35	81	ND	ND		116
	29-Sep-12	10	8.9	270 D	740 D	6.7	ND		1035.6
	13-Nov-13	ND	ND	180	490	ND	ND		670
	20-Feb-15	ND	ND	200	450	ND	ND		650
	31-May-16	ND	ND	92	230	ND	ND		322
	24-Aug-17	3.2	2.9	130	430	3.4	ND		570
	20-Nov-18	ND	ND	84	250	ND	ND		334
	13-Jan-20 24-May-21	ND ND	0.62 1	42 58	110 200	1.2 1.7	ND ND		154 261
	1-Sep-22	4 J	2.4	140	500	5.3	ND ND	3.4	655.1
	7-Jan-91	NA	NA	NA	NA	NA	NA	NA	NA NA
	25-Feb-91	NA	NA	NA	NA	NA	NA	NA	NA
	14-Jul-94	390 J		1100	7400	160 J			9050
	2-Nov-94	100 J		830	4000	61 J			4991
	14-Apr-95	200 J	10	680	3800	130 J			4820
	23-Aug-95	660	<u> </u>	1500	7700	140 J			10000
	27-Oct-99	250 254		1920 J	3510 4320				3760
	8-Feb-00 27-Apr-00	450		1920 J	6430	125			6494 7005
	25-Jul-00	729			12200	162			13091
	19-Oct-00	503		2810	9840	217			13370
	21-Dec-00	197		1670	3240	46.6			5154
MW-201D	28-Feb-01	267		1960	4780				7007
IVIVV-201D	19-Apr-01	252		2300	4220	110			6882
	25-Oct-01	301		2840	4770				7911
	11-Apr-02	103		2450	1850				4403
	29-Oct-02	312		2690	5810	136			8948
	29-Apr-03	277		3030	3980				7287
	27-Oct-03 29-Apr-04	354 201		2890 2620	8430 1890				11674 4711
	28-Oct-04	271		3320	5230	141			8962
	12-Feb-07	190	38	1000	1600	130	ND		2958
	15-Aug-07	2700 D	660	9600 D	46000 D	440	ND		59400
	13-Mar-08	92	21 J	810	3300	40 J	ND		4263
	20-Nov-08	190	34 J	2000	5900	56 J			8180
	4-Feb-10	ND	ND	800	3100	ND	ND		3900
	1-May-11	150	ND	1100	4100	ND	ND		5350
	29-Sep-12	200	ND	1200	5200 D	ND	ND		6600
	13-Nov-13	ND	ND	710	3400	ND	ND		4110
	20-Feb-15 31-May-16	ND ND	ND ND	410 720	2500 4600	ND ND	ND ND		2910 5320
	24-Aug-17	190	ND ND	1100	5900	110	ND ND		7300
	20-Nov-18	ND	ND ND	430	2300	ND	ND ND		2730
	13-Jan-20	49	19	510	2600	44	ND		3222
	24-May-21	150	60	750	5300	64	ND		6324
	1-Sep-22	160	70	1200	6000	54	ND	ND	7484
	7-Jan-91 25-Feb-91	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA		NA NA
	25-Feb-91 14-Jul-94	NA NA	NA NA	NA NA	NA NA	NA NA	NA NA		NA NA
	2-Nov-94	NA NA	NA NA	NA NA	NA NA	NA	NA NA		NA NA
	14-Apr-95	6 J		6 J	42	1 J			55
	23-Aug-95	ļ. — . — <u>.</u> — . — .	2 J	3 J	160	4 J	<u> </u>	 -	169
	27-Oct-99 27-Apr-00	3 3.37			20 33.9		2		25 37
	27-Apr-00 25-Jul-00	3.37 NS	NS	NS	NS NS	NS	NS		NS
SUPPLY	19-Oct-00	186	29.9	44.4	1490				1750
	21-Dec-00	4.3		5.44	52.5				62
	28-Feb-01	6.36		4.68	70				81
	19-Apr-01 25-Oct-01	43.5	5.13	23.4	17.4 456				17 528
	11-Apr-02	3.73	0.13	23.4 5.15	48.5				528 57
			•	•			•	•	•



SUMMARY OF HISTORIC ON-SITE GROUNDWATER ANALYTICAL RESULTS

Enarc-O Machine Products, Inc. Lima, New York NYSDEC Registry No. 8-26-011

WELL	DATE	COMPOUND								
******		1,1,1-TCA	1,1-DCE	cis-1,2-DCE	TCE	PCE	Toluene	1,1-DCA	VOCs	
	29-Oct-02	100	12.2	35.6	980	10.3			1138	
	29-Apr-03	2.94		10.9	47				61	
	27-Oct-03	126	20.4	52.9	1890				2089	
	29-Apr-04				20.5				21	
	28-Oct-04	22.4	2.91	15.7	245	2.1			288	
	12-Feb-07	8.8		11	120				140	
	15-Aug-07	0.91 J		3.1	18				22	
SUPPLY	12-Mar-08	8.1	2	30	180 D	2.3			222	
	20-Nov-08	1.1	2.9	21	240	2.2 J			267	
	4-Feb-10	ND	ND	12	87	ND	ND		99	
	1-May-11	ND	ND	ND	7.9	ND	ND		8	
	29-Sep-12	ND	ND	ND	8.7	ND	ND		9	
	13-Nov-13	ND	ND	5.3	93	ND	ND		98	
	20-Feb-15	ND	ND	ND	15	ND	ND		15	
	31-May-16	ND	ND	ND	9.8	ND	ND		10	
	24-Aug-17	5	3.6	6.2	100	1.8 J	ND		117	
	20-Nov-18	6	ND	12	180	ND	ND		198	
	13-Jan-20	4.9	2.8	ND	180	2.8 J	ND		191	
	24-May-21	ND	0.38 J	2.2 J	18	0.35 J	ND		21	
	1-Sep-22	ND	0.27 J	1.1 J	12	ND	ND	ND	13.37	

- Notes:
 1. All concentrations in ug/L or parts-per-billion (ppb).
 - 2. J = Indicates an estimated concentration.

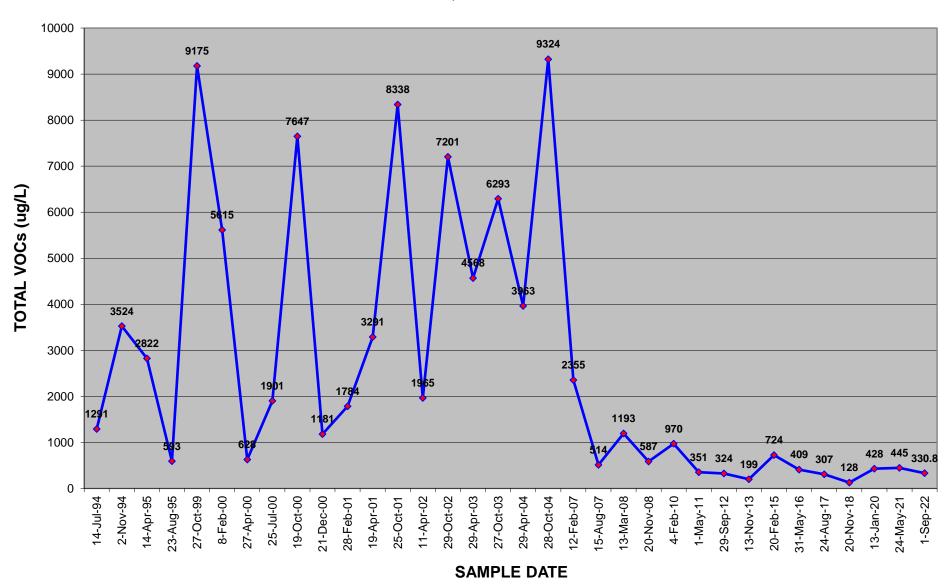
 - U = Indicates compound analyzed for but not detected.
 D = Compound identified at the secondary dillution factor.
 - 5. NA = Not analyzed.

 - NS = Not Sampled.
 ND = None detected (blank space also indicates not detected).
 - 8. Heavy dashed and dotted line indicates time after which LNAPL was observed in MW-201D.
 - 9. Historic concentration data provided by Kadis Enarc-O (pre-2007)
 - 10. Highlighted concentrations indicate the September 2022 sampling event.



HISTORIC ANALYTICAL RESULTS MW-3

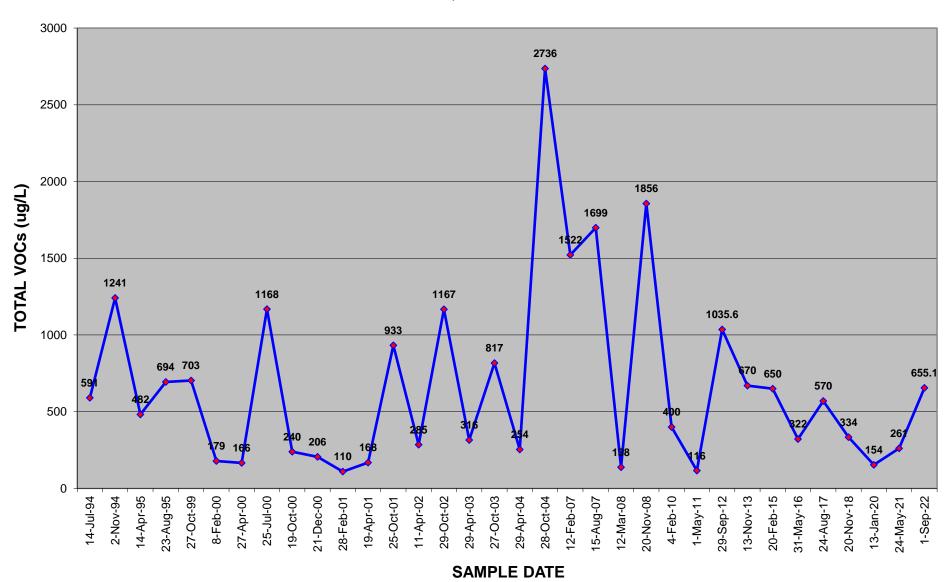
Enarc-O Machine Products Lima, New York





HISTORIC ANALYTICAL RESULTS MW-5

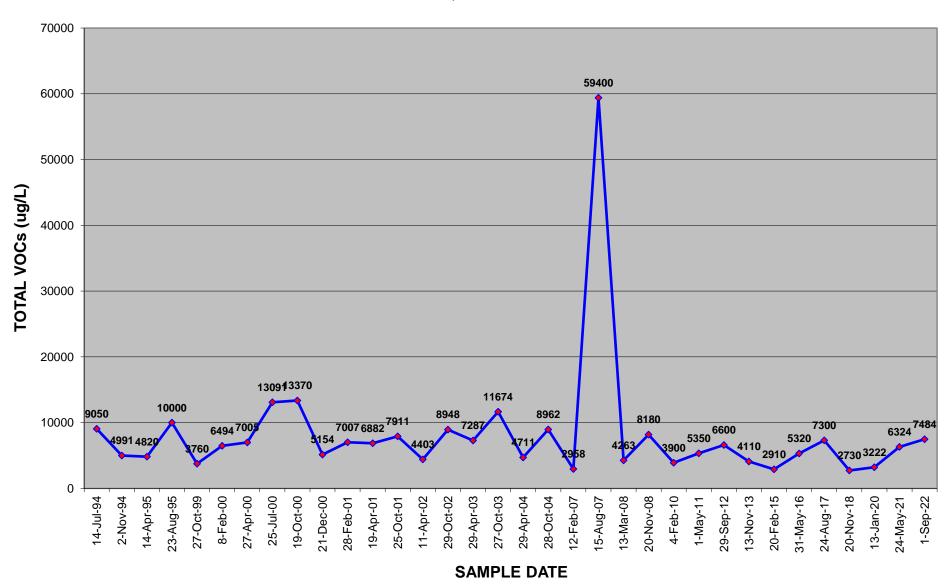
Enarc-O Machine Products Lima, New York





HISTORIC ANALYTICAL RESULTS MW-201D

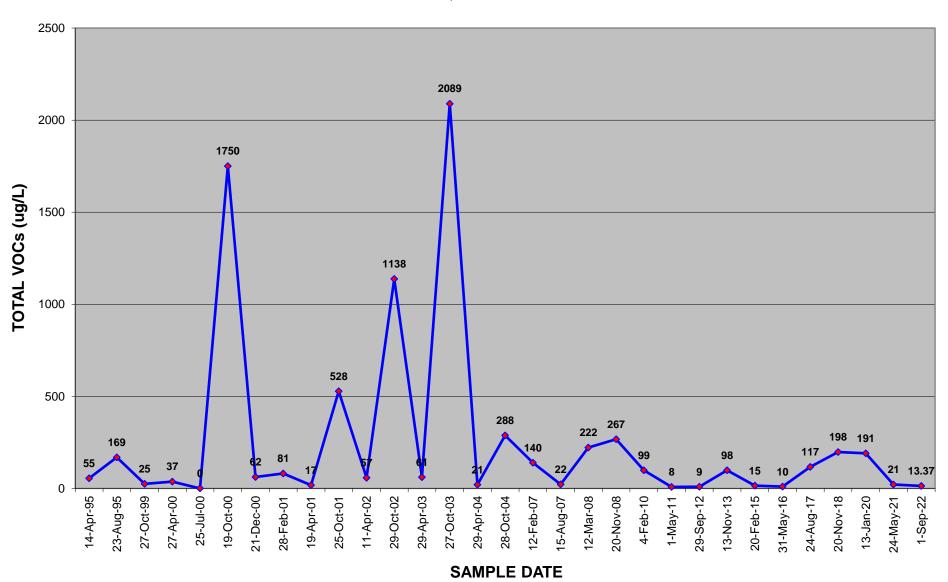
Enarc-O Machine Products Lima, New York





HISTORIC ANALYTICAL RESULTS SUPPLY WELL

Enarc-O Machine Products Lima, New York



WELL REPAIR PHOTOS



PHOTOGRAPHIC LOG

Client Name:Site Location:Project No.:Alco Mfg Corp1175 Bragg St., Honeoye FallsB0672-022-002

Photo No. Date
1 11/04/22

Direction Photo Taken: NW

Description: MW-201D Roadbox Repair



Photo No. 2 11/04/22

Direction Photo Taken:
E

Description:
MWV-202 Roadbox Repair

Prepared By: THF