

#### **DATA PACKAGE**

GENERAL CHEMISTRY
METALS
SEMI-VOLATILE ORGANICS
VOLATILE ORGANICS

**PROJECT NAME: PVSC MONTHLY 2025** 

**ARDMORE CHEMICAL** 

**29 Riverside Avenue** 

Newark, NJ - 07104-

Phone No: 973-481-2406

ORDER ID: Q2264

**ATTENTION: Michael Sharphouse** 







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### **Cover Page**

Order ID:	Q2264
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**Project ID:** PVSC Monthly 2025

**Client:** Ardmore Chemical

Lab Sample Number Client Sample Number

Q2264-01 EFF-WW Q2264-04 EF-WW

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. Release of the data contained in this hard copy data package has been authorized by the laboratory manager or his designee, as verified by the following signature.

NYDOH CERTIFICATION NO - 11376 NJDEP CERTIFICATION NO - 20012

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#### **CASE NARRATIVE**

**Ardmore Chemical** 

**Project Name: PVSC Monthly 2025** 

Project # N/A Order ID # Q2264 Test Name: VOC-PP

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 06/06/2025.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: BOD5, Cyanide, Mercury, Metals Group3, SVOCMS Group1, TSS and VOC-PP. This data package contains results for VOC-PP.

#### C. Analytical Techniques:

The analysis performed on instrument MSVOA\_X were done using GC column DB-624UI 20m 0.18mm 1.0 um. Cat#121-1324UIThe analysis of VOC-PP was based on method 624.1.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria.

The Internal Standards Areas met the acceptable requirements.

The Retention Times were acceptable for all samples.

The RPD met criteria.

The Blank Spike met requirements for all samples.

The Blank Spike Duplicate met requirements for all samples.

The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements.

The Continuous Calibration met the requirements.

The Tuning criteria met requirements.

Samples EFF-WW was diluted due to foamy Sample.

#### E. Additional Comments:

As per method, MS/MSD is required to be performed with the sample analysis. However, Lab did not receive sufficient volume to perform the MS/MSD

therefore MS/MSD were not performed for this project. However, Lab has performed LCS/LCSD instead.

Trip Blank was not provided with this set of samples.

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <35% for the Initial

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Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is > 35% for the Initial Calibration curve for SW-846 analysis.

#### **F. Manual Integration Comments:**

Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. The laboratory manager or his designee, as verified by the following signature has authorized release of the data contained in this hard copy data package.

Signature		

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#### **CASE NARRATIVE**

**Ardmore Chemical** 

**Project Name: PVSC Monthly 2025** 

Project # N/A Order ID # Q2264

**Test Name: SVOCMS Group1** 

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 06/06/2025.

#### **B.** Parameters

According to the Chain of Custody document, the following analyses were requested: BOD5, Cyanide, Mercury, Metals Group3, SVOCMS Group1, TSS and VOC-PP. This data package contains results for SVOCMS Group1.

#### C. Analytical Techniques:

The samples were analyzed on instrument BNA\_F using GC Column DB-UI 8270D which is 20 meters, 0.18 mm ID, 0.36 um dfThe analysis of SVOCMS Group1 was based on method 625.1 and extraction was done based on method 3510.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Surrogate recoveries met the acceptable criteria except for EF-WW [2,4,6-Tribromophenol - 2%, 2-Fluorophenol - 1%, Phenol-d6 - 0% and Terphenyl-d14 - 52%]. Due to matrix interference, which can be observed by the abnormal chromatogram. Reanalyzing this sample will give the same result. Hence this analysis reported as final results.

The Internal Standards Areas met the acceptable requirements except for EF-WW. Due to matrix interference, which can be observed by the abnormal chromatogram. Reanalyzing this sample will give the same result. Hence this analysis reported as final results. The Retention Times were acceptable for all samples.

The RPD met criteria.

The Blank Spike met requirements for all samples.

The Blank Spike Duplicate met requirements for all samples.

The Blank analysis did not indicate the presence of lab contamination.

The Initial Calibration met the requirements.

The Continuous Calibration met the requirements.

The Tuning criteria met requirements.

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#### E. Additional Comments:

Alliance has analyzed samples for SVOCMS Group1 by Method 625.1 for Project "PVSC Monthly 2025". Alliance certification was in applied status for compound "2,4-Dimethylphenol" with NJDEP for Method 625.1 for SVOC group 1 at the time when samples for Project "PVSC Monthly 2025 "were analyzed

As per method, MS/MSD is required to be performed with the sample analysis. However, Lab did not receive sufficient volume to perform the MS/MSD therefore MS/MSD were not performed for this project. However, Lab has performed LCS/LCSD instead.

Please use %D calculated based on Avg RF and CCRF for all compounds using Average Response Factor when the %RSD value for a compound is <35% for the Initial Calibration curve and use %D calculated based on Amount added and Calculated amount for all compounds using Linear Regression when the %RSD value for a compound is > 35% for the Initial Calibration curve for SW-846 analysis.

#### **F. Manual Integration Comments:**

Please refer to the Manual integration Report included with the Run Logs for information on the manual integrations performed.

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Signature		

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#### **CASE NARRATIVE**

**Ardmore Chemical** 

**Project Name: PVSC Monthly 2025** 

Project # N/A **Order ID # Q2264** 

**Test Name: Mercury, Metals Group3** 

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 06/06/2025.

#### **B. Parameters:**

According to the Chain of Custody document, the following analyses were requested: BOD5, Cyanide, Mercury, Metals Group3, SVOCMS Group1, TSS and VOC-PP. This data package contains results for Mercury, Metals Group3.

#### C. Analytical Techniques:

The analysis and digestion of Metals Group3 was based on 200.7 and The analysis and digestion of Mercury was based on 245.1.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Blank Spike met requirements for all samples.

The Duplicate analysis met criteria for all samples.

The Matrix Spike analysis met criteria for all samples.

The Matrix Spike Duplicate analysis met criteria for all samples.

The Blank analysis did not indicate the presence of lab contamination.

The Calibration met the requirements.

The Serial Dilution met the acceptable requirements.

#### E. Additional Comments:

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Signature		
Signature		

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#### CASE NARRATIVE

**Ardmore Chemical** 

**Project Name: PVSC Monthly 2025** 

Project # N/A Order ID # Q2264

Test Name: BOD5, Cyanide, TSS

#### A. Number of Samples and Date of Receipt:

2 Water samples were received on 06/06/2025.

#### **B. Parameters:**

According to the Chain of Custody document, the following analyses were requested: BOD5, Cyanide, Mercury, Metals Group3, SVOCMS Group1, TSS and VOC-PP. This data package contains results for BOD5, Cyanide, TSS.

#### C. Analytical Techniques:

The analysis of TSS was based on method SM2540 D, The analysis of Cyanide was based on method SM4500-CN C,E and The analysis of BOD5 was based on method SM5210 B.

#### D. QA/ QC Samples:

The Holding Times were met for all analysis.

The Blank Spike met requirements for all samples.

The Duplicate analysis met criteria for all samples.

The Matrix Spike analysis met criteria for all samples.

The Matrix Spike Duplicate analysis met criteria for all samples.

The Blank analysis did not indicate the presence of lab contamination.

The Calibration met the requirements.

#### **E. Additional Comments:**

I certify that the data package is in compliance with the terms and conditions of the contract, both technically and for completeness, for other than the conditions detailed above. The laboratory manager or his designee, as verified by the following signature has authorized release of the data contained in this hard copy data package.

Signature		
Signature		

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#### DATA REPORTING QUALIFIERS- INORGANIC

For reporting results, the following "Results Qualifiers" are used:

- J Indicates the reported value was obtained from a reading that was less than the Contract Required Detection Limit (CRDL), but greater than or equal to the Instrument Detection Limit (IDL).
- U Indicates the analyte was analyzed for, but not detected.
- ND Indicates the analyte was analyzed for, but not detected
- E Indicates the reported value is estimated because of the presence of interference
- M Indicates Duplicate injection precision not met.
- N Indicates the spiked sample recovery is not within control limits.
- S Indicates the reported value was determined by the Method of Standard Addition (MSA).
- \* Indicates that the duplicate analysis is not within control limits.
- + Indicates the correlation coefficient for the MSA is less than 0.995.
- D Indicates the reported value is from a secondary analysis with a dilution factor. The original analysis exceeded the calibration range.
- M Method qualifiers
  - **"P"** for ICP instrument
  - "PM" for ICP when Microwave Digestion is used
  - "CV" for Manual Cold Vapor AA
  - "AV" for automated Cold Vapor AA
  - "CA" for MIDI-Distillation Spectrophotometric "AS" for Semi –Automated Spectrophotometric
  - "C" for Manual Spectrophotometric
  - **"T"** for Titrimetric
  - "NR" for analyte not required to be analyzed
- OR Indicates the analyte's concentration exceeds the calibrated range of the
  - instrument for that specific analysis.
- Q Indicates the LCS did not meet the control limits requirements
- H Sample Analysis Out Of Hold Time



#### DATA REPORTING QUALIFIERS- ORGANIC

For reporting results, the following "Results Qualifiers" are used:

Value	If the result is a value greater than or equal to the detection limit, report the value
U	Indicates the compound was analyzed for but was not detected. Report the minimum detection limit for the sample with the U, i.e. "10 U". This is not necessarily the instrument detection limit attainable for this particular sample based on any concentration or dilution that may have been required.
ND	Indicates the analyte was analyzed for, but not detected
<b>J</b>	<ul> <li>Indicates an estimated value. This flag is used:</li> <li>(1) When estimating a concentration for a tentatively identified compound (library search hits, where a 1:1 response is assumed.)</li> <li>(2) When the mass spectral data indicated the identification, however the result was less than the specified detection limit greater than zero. If the detection limit was 10ug/L and a concentration of 3 ug/L was calculated report as 3 J. This is flag is used when similar situation arise on any organic parameter i.e. Pest, PCB and others.</li> <li>Indicates the analyte was found in the blank as well as the sample report as</li> </ul>
	"12 B".
E	Indicates the analyte's concentration exceeds the calibrated range of the instrument for that specific analysis.
D	This flag identifies all compounds identified in an analysis at a secondary dilution factor.
P	This flag is used for Pesticide/PCB target analyte when there is $>25\%$ difference for detected concentrations between the two GC columns. The lower of the two values is reported on Form 1 and flagged with a "P".
N	This flag indicates presumptive evidence of a compound. This is only used for tentatively identified compounds (TICs), where the identification is based on a mass spectral library search. It applies to all TIC results. For generic characterization of a TIC, such as chlorinated hydrocarbon, the flag is not used.
A	This flag indicates that a Tentatively Identified Compound is a suspected aldol-condensation product.
Q	Indicates the LCS did not meet the control limits requirements

Aliance

#### APPENDIX A

#### **QA REVIEW GENERAL DOCUMENTATION**

Project #: Q2264

	Completed
For thorough review, the report must have the following:	
GENERAL:	
Are all original paperwork present (chain of custody, record of communication, airbill, sample management lab chronicle, login page)	<u> </u>
Check chain-of-custody for proper relinquish/return of samples	<u> </u>
Is the chain of custody signed and complete	<u> </u>
Check internal chain-of-custody for proper relinquish/return of samples /sample extracts	<del>'</del> <del>'</del> <del>'</del> <del>'</del> <del>'</del> <del>'</del>
Collect information for each project id from server. Were all requirements followed	<u> </u>
COVER PAGE:	
Do numbers of samples correspond to the number of samples in the Chain of Custody on login page	<u> </u>
Do lab numbers and client Ids on cover page agree with the Chain of Custody	<u> </u>
CHAIN OF CUSTODY:	
Do requested analyses on Chain of Custody agree with form I results	<u> </u>
Do requested analyses on Chain of Custody agree with the log-in page	<u> </u>
Were the correct method log-in for analysis according to the Analytical Request and Chain of Castody	<u> </u>
Were the samples received within hold time	<u> </u>
Were any problems found with the samples at arrival recorded in the Sample Management Laboratory Chronicle	<u> </u>
ANALYTICAL:	
Was method requirement followed?	<u> </u>
Was client requirement followed?	<u> </u>
Does the case narrative summarize all QC failure?	<del>'</del> <del>'</del> <del>'</del> <del>'</del> <del>'</del> <del>'</del>
All runlogs and manual integration are reviewed for requirements	<u> </u>
All manual calculations and /or hand notations verified	<u> </u>

QA Review Signature: SOHIL JODHANI Date: 06/20/2025

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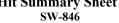


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## **Hit Summary Sheet**

SDG No.: Q2264

Client: Ardmore Chemical





Sample ID	Client ID	Matrix	Parameter	Concentration	C MDL	RDL	Units
Client ID: Q2264-01	EFF-WW EFF-WW	Water	Chloroform	12.0	J 2.80	25.0	ug/L
			Total Voc:	12.0			
			<b>Total Concentration:</b>	12.0			

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# SAMPLE DATA

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LOW



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Level:

#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 Project: PVSC Monthly 2025 Date Received: 06/06/25 Client Sample ID: EFF-WW SDG No.: Q2264 Lab Sample ID: Matrix: Water Q2264-01 Analytical Method: E624.1 % Solid: Sample Wt/Vol: 5 Final Vol: 5000 Units: mLSoil Aliquot Vol: uL Test: VOC-PP

GC Column:

Prep Method:

File ID/Qc Batch: Dilution: Prep Date Date Analyzed Prep Batch ID VX046634.D 5 06/11/25 13:40 VX061125

ID: 0.18

DB-624UI

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
TARGETS						
74-87-3	Chloromethane	3.20	U	3.20	25.0	ug/L
75-01-4	Vinyl Chloride	4.20	U	4.20	25.0	ug/L
74-83-9	Bromomethane	4.00	U	4.00	25.0	ug/L
75-00-3	Chloroethane	11.6	U	11.6	25.0	ug/L
75-69-4	Trichlorofluoromethane	4.00	U	4.00	25.0	ug/L
75-35-4	1,1-Dichloroethene	3.80	U	3.80	25.0	ug/L
107-02-8	Acrolein	33.1	U	33.1	130	ug/L
107-13-1	Acrylonitrile	14.0	U	14.0	130	ug/L
75-09-2	Methylene Chloride	4.30	U	4.30	25.0	ug/L
156-60-5	trans-1,2-Dichloroethene	4.10	U	4.10	25.0	ug/L
75-34-3	1,1-Dichloroethane	3.40	U	3.40	25.0	ug/L
56-23-5	Carbon Tetrachloride	3.70	U	3.70	25.0	ug/L
67-66-3	Chloroform	12.0	J	2.80	25.0	ug/L
71-55-6	1,1,1-Trichloroethane	3.20	U	3.20	25.0	ug/L
71-43-2	Benzene	2.30	U	2.30	25.0	ug/L
107-06-2	1,2-Dichloroethane	2.50	U	2.50	25.0	ug/L
79-01-6	Trichloroethene	2.50	U	2.50	25.0	ug/L
78-87-5	1,2-Dichloropropane	2.30	U	2.30	25.0	ug/L
75-27-4	Bromodichloromethane	3.20	U	3.20	25.0	ug/L
108-88-3	Toluene	2.30	U	2.30	25.0	ug/L
10061-02-6	t-1,3-Dichloropropene	3.60	U	3.60	25.0	ug/L
10061-01-5	cis-1,3-Dichloropropene	3.40	U	3.40	25.0	ug/L
79-00-5	1,1,2-Trichloroethane	2.30	U	2.30	25.0	ug/L
110-75-8	2-Chloroethyl vinyl ether	23.2	U	23.2	130	ug/L
124-48-1	Dibromochloromethane	3.30	U	3.30	25.0	ug/L
127-18-4	Tetrachloroethene	4.20	U	4.20	25.0	ug/L
108-90-7	Chlorobenzene	2.40	U	2.40	25.0	ug/L
100-41-4	Ethyl Benzene	2.80	U	2.80	25.0	ug/L
179601-23-1	m/p-Xylenes	6.50	U	6.50	50.0	ug/L
95-47-6	o-Xylene	3.40	U	3.40	25.0	ug/L

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Level:

LOW

#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 Date Received: Project: PVSC Monthly 2025 06/06/25 SDG No.: Q2264 Client Sample ID: EFF-WW Lab Sample ID: Q2264-01 Matrix: Water Analytical Method: E624.1 % Solid: Final Vol: Sample Wt/Vol: 5 Units: mL5000 uL Soil Aliquot Vol: Test: VOC-PP uL

GC Column: DB-624UI ID: 0.18

Prep Method:

File ID/Qc Batch: Dilution: Prep Date Date Analyzed Prep Batch ID VX046634.D 5 06/11/25 13:40 VX061125

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
75-25-2	Bromoform	4.70	U	4.70	25.0	ug/L
79-34-5	1,1,2,2-Tetrachloroethane	2.20	U	2.20	25.0	ug/L
541-73-1	1,3-Dichlorobenzene	3.40	U	3.40	25.0	ug/L
106-46-7	1,4-Dichlorobenzene	4.10	U	4.10	25.0	ug/L
95-50-1	1,2-Dichlorobenzene	3.40	U	3.40	25.0	ug/L
SURROGATES						
17060-07-0	1,2-Dichloroethane-d4	32.5		91 - 110	108%	SPK: 30
2037-26-5	Toluene-d8	29.0		91 - 112	97%	SPK: 30
460-00-4	4-Bromofluorobenzene	29.5		63 - 112	98%	SPK: 30
INTERNAL ST	ANDARDS					
74-97-5	Bromochloromethane	16600	4.922			
540-36-3	1,4-Difluorobenzene	91500	6.769			
3114-55-4	Chlorobenzene-d5	85900	10.055			

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

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#### LAB CHRONICLE

**OrderID:** Q2264 **OrderDate:** 6/6/2025 2:07:00 PM

Client:Ardmore ChemicalProject:PVSC Monthly 2025Contact:Michael SharphouseLocation:D41,VOA Ref. #3 Water

ClientID Sample Date **Prep Date** Received LabID Matrix Test Method **Anal Date** Q2264-01 06/06/25 06/06/25 **EFF-WW** Water VOC-PP 624.1 06/11/25

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Α

В

C



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#### Hit Summary Sheet SW-846

**SDG No.:** Q2264

Client: Ardmore Chemical

Sample ID Client ID Matrix Parameter Concentration C MDL RDL Units

Client ID:

0.000

Total Svoc: 0.00
Total Concentration: 0.00

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# SAMPLE DATA

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#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 Project: PVSC Monthly 2025 Date Received: 06/06/25 Client Sample ID: EF-WW SDG No.: Q2264 Lab Sample ID: Q2264-04 Matrix: Water % Solid: 0 Analytical Method: 625.1 Sample Wt/Vol: 870 Units: mL Final Vol: 1000 uL SVOCMS Group1 Soil Aliquot Vol: uL Test:

Extraction Type: Decanted: N Level: LOW

Injection Volume : GPC Factor : 1.0 GPC Cleanup : N PH :

Prep Method: 3510C

 File ID/Qc Batch:
 Dilution:
 Prep Date
 Date Analyzed
 Prep Batch ID

 BF142732.D
 1
 06/10/25 08:46
 06/11/25 13:52
 PB168378

B1112702.B	•	00/10/20	000	00/11/20 13.02					
CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units			
TARGETS									
62-75-9	n-Nitrosodimethylamine	0.99	U	0.99	11.5	ug/L			
108-95-2	Phenol	1.00	U	1.00	5.70	ug/L			
111-44-4	bis(2-Chloroethyl)ether	0.93	U	0.93	5.70	ug/L			
95-57-8	2-Chlorophenol	0.67	U	0.67	5.70	ug/L			
108-60-1	2,2-oxybis(1-Chloropropane)	1.50	U	1.50	5.70	ug/L			
621-64-7	n-Nitroso-di-n-propylamine	1.60	U	1.60	5.70	ug/L			
67-72-1	Hexachloroethane	0.75	U	0.75	5.70	ug/L			
98-95-3	Nitrobenzene	0.87	U	0.87	5.70	ug/L			
78-59-1	Isophorone	0.86	U	0.86	5.70	ug/L			
88-75-5	2-Nitrophenol	2.00	U	2.00	5.70	ug/L			
105-67-9	2,4-Dimethylphenol	2.10	U	2.10	5.70	ug/L			
111-91-1	bis(2-Chloroethoxy)methane	0.78	U	0.78	5.70	ug/L			
120-83-2	2,4-Dichlorophenol	0.60	U	0.60	5.70	ug/L			
120-82-1	1,2,4-Trichlorobenzene	0.62	U	0.62	5.70	ug/L			
91-20-3	Naphthalene	0.57	U	0.57	5.70	ug/L			
87-68-3	Hexachlorobutadiene	0.62	U	0.62	5.70	ug/L			
59-50-7	4-Chloro-3-methylphenol	0.68	U	0.68	5.70	ug/L			
77-47-4	Hexachlorocyclopentadiene	4.20	U	4.20	11.5	ug/L			
88-06-2	2,4,6-Trichlorophenol	0.59	U	0.59	5.70	ug/L			
91-58-7	2-Chloronaphthalene	0.70	U	0.70	5.70	ug/L			
131-11-3	Dimethylphthalate	0.70	U	0.70	5.70	ug/L			
208-96-8	Acenaphthylene	0.86	U	0.86	5.70	ug/L			
606-20-2	2,6-Dinitrotoluene	1.10	U	1.10	5.70	ug/L			
83-32-9	Acenaphthene	0.63	U	0.63	5.70	ug/L			
51-28-5	2,4-Dinitrophenol	6.90	U	6.90	11.5	ug/L			
100-02-7	4-Nitrophenol	2.70	U	2.70	11.5	ug/L			
121-14-2	2,4-Dinitrotoluene	1.40	U	1.40	5.70	ug/L			
84-66-2	Diethylphthalate	0.79	U	0.79	5.70	ug/L			
7005-72-3	4-Chlorophenyl-phenylether	0.78	U	0.78	5.70	ug/L			

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#### Report of Analysis

Client: Ardmore Chemical Date Collected: 06/06/25

Project: PVSC Monthly 2025 Date Received: 06/06/25

Client Sample ID: EF-WW SDG No.: Q2264

Lab Sample ID: Q2264-04 Matrix: Water
Analytical Method: 625.1 % Solid: 0

Sample Wt/Vol: 870 Units: mL Final Vol: 1000 uL
Soil Aliquot Vol: uL Test: SVOCMS Group1

Extraction Type: Decanted: N Level: LOW

Injection Volume : GPC Factor : 1.0 GPC Cleanup : N PH :

Prep Method: 3510C

 File ID/Qc Batch:
 Dilution:
 Prep Date
 Date Analyzed
 Prep Batch ID

 BF142732.D
 1
 06/10/25 08:46
 06/11/25 13:52
 PB168378

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units
86-73-7	Fluorene	0.72	U	0.72	5.70	ug/L
534-52-1	4,6-Dinitro-2-methylphenol	3.30	U	3.30	11.5	ug/L
86-30-6	n-Nitrosodiphenylamine	0.67	U	0.67	5.70	ug/L
103-33-3	Azobenzene	0.93	U	0.93	5.70	ug/L
101-55-3	4-Bromophenyl-phenylether	0.46	U	0.46	5.70	ug/L
118-74-1	Hexachlorobenzene	0.60	U	0.60	5.70	ug/L
87-86-5	Pentachlorophenol	1.80	U	1.80	11.5	ug/L
85-01-8	Phenanthrene	0.57	U	0.57	5.70	ug/L
120-12-7	Anthracene	0.70	U	0.70	5.70	ug/L
84-74-2	Di-n-butylphthalate	1.40	U	1.40	5.70	ug/L
206-44-0	Fluoranthene	0.94	U	0.94	5.70	ug/L
92-87-5	Benzidine	4.90	U	4.90	11.5	ug/L
129-00-0	Pyrene	0.57	U	0.57	5.70	ug/L
85-68-7	Butylbenzylphthalate	2.20	U	2.20	5.70	ug/L
91-94-1	3,3-Dichlorobenzidine	1.10	U	1.10	11.5	ug/L
56-55-3	Benzo(a)anthracene	0.52	U	0.52	5.70	ug/L
218-01-9	Chrysene	0.51	U	0.51	5.70	ug/L
117-81-7	Bis(2-ethylhexyl)phthalate	1.80	U	1.80	5.70	ug/L
117-84-0	Di-n-octyl phthalate	2.70	U	2.70	11.5	ug/L
205-99-2	Benzo(b)fluoranthene	0.56	U	0.56	5.70	ug/L
207-08-9	Benzo(k)fluoranthene	0.55	U	0.55	5.70	ug/L
50-32-8	Benzo(a)pyrene	0.63	U	0.63	5.70	ug/L
193-39-5	Indeno(1,2,3-cd)pyrene	0.68	U	0.68	5.70	ug/L
53-70-3	Dibenzo(a,h)anthracene	0.77	U	0.77	5.70	ug/L
191-24-2	Benzo(g,h,i)perylene	0.79	U	0.79	5.70	ug/L
SURROGATES						
367-12-4	2-Fluorophenol	0.79	*	60 - 140	1%	SPK: 100
13127-88-3	Phenol-d6	0	*	60 - 140	0%	SPK: 100
4165-60-0	Nitrobenzene-d5	67.8		60 - 140	68%	SPK: 100
321-60-8	2-Fluorobiphenyl	60.0		60 - 140	60%	SPK: 100

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#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 Project: PVSC Monthly 2025 Date Received: 06/06/25

Client Sample ID: EF-WW SDG No.: Q2264

Lab Sample ID: Q2264-04 Matrix: Water Analytical Method: 625.1 % Solid: 0

Sample Wt/Vol: 870 Units: Final Vol: 1000 uL mL

Test:

SVOCMS Group1 Soil Aliquot Vol: Level: Extraction Type: Decanted: Ν LOW

GPC Cleanup: Injection Volume: GPC Factor: 1.0 Ν PH:

3510C Prep Method:

File ID/Qc Batch: Dilution: Prep Date Prep Batch ID Date Analyzed BF142732.D 1 06/10/25 08:46 06/11/25 13:52 PB168378

CAS Number	Parameter	Conc.	Qualifier	MDL	LOQ / CRQL	Units	
118-79-6	2,4,6-Tribromophenol	2.09	*	60 - 140	2%	SPK: 100	
1718-51-0	Terphenyl-d14	51.6	*	60 - 140	52%	SPK: 100	
INTERNAL STA	ANDARDS						
3855-82-1	1,4-Dichlorobenzene-d4	50300	6.892				
1146-65-2	Naphthalene-d8	171000	8.186				
15067-26-2	Acenaphthene-d10	83200	9.933				
1517-22-2	Phenanthrene-d10	136000	11.422				
1719-03-5	Chrysene-d12	137000	14.068				
1520-96-3	Perylene-d12	151000	15.568				

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

E = Value Exceeds Calibration Range

Q = indicates LCS control criteria did not meet requirements

M = MS/MSD acceptance criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

N = Presumptive Evidence of a Compound

\* = Values outside of QC limits

D = Dilution

() = Laboratory InHouse Limit

A = Aldol-Condensation Reaction Products

Q2264

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#### LAB CHRONICLE

**OrderID:** Q2264 **OrderDate:** 6/6/2025 2:07:00 PM

Client:Ardmore ChemicalProject:PVSC Monthly 2025Contact:Michael SharphouseLocation:D41,VOA Ref. #3 Water

Sample Date **Prep Date** Received LabID ClientID Matrix Test Method **Anal Date** Q2264-04 06/06/25 06/06/25 EF-WW Water SVOCMS Group1 625.1 06/10/25 06/11/25

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Q2264

SDG No.:

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#### **Hit Summary Sheet** SW-846

Q2264 Order ID:

Client:	Ardmore Chemical			Project ID	:	PVSC Monthly 2025		
Sample ID	Client ID	Matrix	Parameter	Concentration	C	MDL	RDL	Units
Client ID:	EF-WW							
Q2264-04	EF-WW	Water	Lead	21.8		1.21	6.00	ug/L
Q2264-04	EF-WW	Water	Mercury	0.039	J	0.027	0.20	ug/L
Q2264-04	EF-WW	Water	Zinc	237		1.60	20.0	ug/L

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# SAMPLE DATA

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D



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#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 Project: PVSC Monthly 2025 Date Received: 06/06/25 Client Sample ID: EF-WW SDG No.: Q2264 Lab Sample ID: Q2264-04 Matrix: Water % Solid: Level (low/med): 0 low

Cas	Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Prep Date	Date Ana.	Ana Met.	Prep Met.
7439-92-1	Lead	21.8		1	1.21	6.00	ug/L	06/12/25 12:15	06/17/25 13:53	EPA 200.7	E200.7
7439-97-6	Mercury	0.039	J	1	0.027	0.20	ug/L	06/11/25 08:30	06/11/25 10:11	E245.1	
7440-66-6	Zinc	237		1	1.60	20.0	ug/L	06/12/25 12:15	06/17/25 13:53	EPA 200.7	E200.7

Color Before: Colorless

Clarity Before: Clarity After: Clear Clear Texture:
Artifacts:

Color After: Colorless

Mercury

U = Not Detected

Comments:

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

D = Dilution

Q = indicates LCS control criteria did not meet requirements

J = Estimated Value

B = Analyte Found in Associated Method Blank

\* = indicates the duplicate analysis is not within control limits.

E = Indicates the reported value is estimated because of the presence

of interference.

OR = Over Range

N =Spiked sample recovery not within control limits

Q2264



#### LAB CHRONICLE

Q2264 OrderID: 6/6/2025 2:07:00 PM OrderDate:

Ardmore Chemical Client: PVSC Monthly 2025 Project: Contact:

Michael Sharphouse Location: D41,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q2264-04	EF-WW	Water			06/06/25			06/06/25
			Mercury	245.1		06/11/25	06/11/25	
			Metals Group3	200.7		06/12/25	06/17/25	

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# SAMPLE DATA











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#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 12:30

Project: PVSC Monthly 2025 Date Received: 06/06/25

Client Sample ID: EFF-WW SDG No.: Q2264

Lab Sample ID: Q2264-01 Matrix: WATER

% Solid: 0

Parameter	Conc.	Qua.	DF	MDL	LOQ / CRQL	Units	Prep Date	Date Ana.	Ana Met.
Cyanide	0.0012	U	1	0.0012	0.0050	mg/L	06/09/25 11:50	06/09/25 14:01	SM 4500-CN
									C-16 plus E-16

Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

D = Dilution

Q = indicates LCS control criteria did not meet requirements

H = Sample Analysis Out Of Hold Time

J = Estimated Value

B = Analyte Found in Associated Method Blank

\* = indicates the duplicate analysis is not within control limits.

E = Indicates the reported value is estimated because of the presence of interference.

OR = Over Range

N =Spiked sample recovery not within control limits

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#### **Report of Analysis**

Client: Ardmore Chemical Date Collected: 06/06/25 12:30

Project: PVSC Monthly 2025 Date Received: 06/06/25

Client Sample ID: EF-WW SDG No.: Q2264

Lab Sample ID: Q2264-04 Matrix: WATER

% Solid: 0

Parameter	Conc. Qua.	DF	MDL	LOQ / CRQL	Units	Prep Date	Date Ana.	Ana Met.
BOD5	739	1	0.20	2.00	mg/L		06/06/25 16:10	SM 5210 B-16
TSS	24.4	1	1.00	4.00	mg/L		06/09/25 13:00	SM 2540 D-15

Comments:

U = Not Detected

LOQ = Limit of Quantitation

MDL = Method Detection Limit

LOD = Limit of Detection

D = Dilution

Q = indicates LCS control criteria did not meet requirements

H = Sample Analysis Out Of Hold Time

J = Estimated Value

B = Analyte Found in Associated Method Blank

\* = indicates the duplicate analysis is not within control limits.

E = Indicates the reported value is estimated because of the presence of interference.

OR = Over Range

N =Spiked sample recovery not within control limits



LAB CHRONICLE

**OrderID:** Q2264 **OrderDate:** 6/6/2025 2:07:00 PM

Client:Ardmore ChemicalProject:PVSC Monthly 2025Contact:Michael SharphouseLocation:D41,VOA Ref. #3 Water

LabID	ClientID	Matrix	Test	Method	Sample Date	Prep Date	Anal Date	Received
Q2264-01	EFF-WW	WATER			06/06/25 12:30			06/06/25
			Cyanide	SM4500-CN C,E		06/09/25	06/09/25 14:01	
Q2264-04	EF-WW	WATER			06/06/25 12:30			06/06/25
			BOD5	SM5210 B			06/06/25	
							16:10	
			TSS	SM2540 D			06/09/25	
							13:00	

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## SHIPPING DOCUMENTS

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ALLIANCE PROJECT NO. QUOTE NO.

COC Number

	CLIENT INFORMATION						CLIENT P	ROJECT IN	FORM/	ATION	TA			CLIENT BILLING INFORMATION						
COMPANY: F	rolm	rttobe sentto: ove Inc	2	PROJE	CT.N	MAN	E:						BILL.	TO:			00		PO#:	
			e Blg#14	PROJEC	CT NO	D.:		LOCA	TION:				ADDF	RESS:						
CITY Neu	NARK	STATE: N	5 ZIP: 07/04	PROJEC	CT MA	ANAG	BER:						CITY			STA	STATE: ZIP:			
		Sharph								ATTE	NTION:				PHC	NE:				
	3 481 240		000	PHONE				FA	V					HE			AN	ALYSIS		- 0
		ROUND INFORMAT	ION	PHONE		DATA	DELIVE	RABLE IN		ATION		0.50.4		اللبر		90		إ	الرسار	
FAX (RUSH) HARDCOPY (D EDD: *TO BE APPRO STANDARD HA	DAYS*	□ Leve	2 (Re 3 (Re w Dat	sults - sults - a)	+ QC) 🗆 + QC 🚨	Level 4 (QC NJ Reduced NYS ASP A Other	d 🗆 US	EPA C				5		1015 7015	/8	/9		MANIFACE		
ALLIANCE SAMPLE ID	s	PROJECT AMPLE IDENTIFICA	ATION	SAMPLE MATRIX	SAM	GRAB BTAI		MPLE ECTION TIME	# OF BOTTLES	1	2	3	4	SERVA 5	6	- 7	8	9		MMENTS fy Preservatives D-NaOH E-ICE F-OTHER
1.	EFF	ww		WW		X	61066	35 12 3		×	X									
2.		WW		WW	X		6/6/25					X	X	X						
3.												-								
4.																				
5.																				
6.																				
7.																				
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3.	NQUISHED BY SAMPLER: DATE/TIME: RECEIVED BY 3.						Page	of		CLIENT	I: Q	Hand D	elivered	0 0	ther					t Complete



#### Laboratory Certification

Certified By	License No.
CAS EPA CLP Contract	68HERH20D0011
5.10 2.7752. 33.111131	001.2.1.1.200011
Connecticut	PH-0830
DOD ELAP (ANAB)	L2219
Maine	2024021
Maryland	296
New Hampshire	255424 Rev 1
New Jersey	20012
New sersey	20012
New York	11376
Pennsylvania	68-00548
Cail Dawrit	505 04 004 00444
Soil Permit	525-24-234-08441
Texas	T104704488

QA Control Code: A2070148

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Fax: 908 789 8922

#### LOGIN REPORT/SAMPLE TRANSFER

Order ID: Q2264

ARDM01

Order Date: 6/6/2025 2:07:00 PM

Project Mgr:

Client Name: Ardmore Chemical

Project Name: PVSC Monthly 2025

Report Type: Level 1

Client Contact: Michael Sharphouse

Receive DateTime: 6/6/2025 1:55:00 PM

**EDD Type:** NONE

Invoice Name: Ardmore Chemical

Purchase Order:

Hard Copy Date:

Invoice Contact: Michael Sharphouse

Date Signoff:

LAB ID	CLIENT ID	MATRIX SAMPLE SAMPLE DATE TIME	TEST	TEST GROUP	METHOD	FAX DATE	DUE DATES
Q2264-01	EFF-WW	Water 06/06/2025 12:30					
			VOC-PP		624.1	10 Bus. Days	

Relinguished By:

Date / Time : 6 10 25

Received By:

Date/Time: Oblob/25 14.55 Rath 5

Storage Area: VOA Refridgerator Room

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