

## Report of Analysis

Client: Ardmore Chemical  
Project: PVSC Monthly 2025  
Client Sample ID: PB170210BL  
Lab Sample ID: PB170210BL  
Analytical Method: 625.1 Level : LOW  
Sample Wt/Vol: 1000 mL Final Vol: 1000 uL  
Prep Method : 3510C Prep Date: 10/22/25

Date Collected:  
Date Received:  
SDG No.: Q3385  
Matrix: Water  
% Solid: 0  
Test: SVOCMS Group1

| CAS Number     | Parameter                   | Conc. | Qua. | DF | MDL  | LOQ / CRQL | Units | Date Ana.      | Prep BatchID |
|----------------|-----------------------------|-------|------|----|------|------------|-------|----------------|--------------|
| <b>TARGETS</b> |                             |       |      |    |      |            |       |                |              |
| 62-75-9        | n-Nitrosodimethylamine      | 0.86  | U    | 1  | 0.86 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 108-95-2       | Phenol                      | 0.91  | U    | 1  | 0.91 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 111-44-4       | bis(2-Chloroethyl)ether     | 0.81  | U    | 1  | 0.81 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 95-57-8        | 2-Chlorophenol              | 0.58  | U    | 1  | 0.58 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 1.30  | U    | 1  | 1.30 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 1.40  | U    | 1  | 1.40 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 67-72-1        | Hexachloroethane            | 0.65  | U    | 1  | 0.65 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 98-95-3        | Nitrobenzene                | 0.76  | U    | 1  | 0.76 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 78-59-1        | Isophorone                  | 0.75  | U    | 1  | 0.75 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 88-75-5        | 2-Nitrophenol               | 1.80  | U    | 1  | 1.80 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 105-67-9       | 2,4-Dimethylphenol          | 1.90  | U    | 1  | 1.90 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 0.68  | U    | 1  | 0.68 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 120-83-2       | 2,4-Dichlorophenol          | 0.52  | U    | 1  | 0.52 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 120-82-1       | 1,2,4-Trichlorobenzene      | 0.54  | U    | 1  | 0.54 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 91-20-3        | Naphthalene                 | 0.50  | U    | 1  | 0.50 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 87-68-3        | Hexachlorobutadiene         | 0.54  | U    | 1  | 0.54 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 59-50-7        | 4-Chloro-3-methylphenol     | 0.59  | U    | 1  | 0.59 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 77-47-4        | Hexachlorocyclopentadiene   | 3.60  | U    | 1  | 3.60 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 88-06-2        | 2,4,6-Trichlorophenol       | 0.51  | U    | 1  | 0.51 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 91-58-7        | 2-Chloronaphthalene         | 0.61  | U    | 1  | 0.61 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 131-11-3       | Dimethylphthalate           | 0.61  | U    | 1  | 0.61 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 208-96-8       | Acenaphthylene              | 0.75  | U    | 1  | 0.75 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 606-20-2       | 2,6-Dinitrotoluene          | 0.92  | U    | 1  | 0.92 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 83-32-9        | Acenaphthene                | 0.55  | U    | 1  | 0.55 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 51-28-5        | 2,4-Dinitrophenol           | 6.00  | U    | 1  | 6.00 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 100-02-7       | 4-Nitrophenol               | 2.40  | U    | 1  | 2.40 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 121-14-2       | 2,4-Dinitrotoluene          | 1.20  | U    | 1  | 1.20 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 84-66-2        | Diethylphthalate            | 0.69  | U    | 1  | 0.69 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 7005-72-3      | 4-Chlorophenyl-phenylether  | 0.68  | U    | 1  | 0.68 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 86-73-7        | Fluorene                    | 0.63  | U    | 1  | 0.63 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 534-52-1       | 4,6-Dinitro-2-methylphenol  | 2.90  | U    | 1  | 2.90 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 86-30-6        | n-Nitrosodiphenylamine      | 0.58  | U    | 1  | 0.58 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 103-33-3       | Azobenzene                  | 0.81  | U    | 1  | 0.81 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 101-55-3       | 4-Bromophenyl-phenylether   | 0.40  | U    | 1  | 0.40 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 118-74-1       | Hexachlorobenzene           | 0.52  | U    | 1  | 0.52 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 87-86-5        | Pentachlorophenol           | 1.60  | U    | 1  | 1.60 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 85-01-8        | Phenanthrene                | 0.50  | U    | 1  | 0.50 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 120-12-7       | Anthracene                  | 0.61  | U    | 1  | 0.61 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 84-74-2        | Di-n-butylphthalate         | 1.20  | U    | 1  | 1.20 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 206-44-0       | Fluoranthene                | 0.82  | U    | 1  | 0.82 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 92-87-5        | Benzidine                   | 4.30  | U    | 1  | 4.30 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 129-00-0       | Pyrene                      | 0.50  | U    | 1  | 0.50 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |

## Report of Analysis

|                    |                   |                     |                 |               |
|--------------------|-------------------|---------------------|-----------------|---------------|
| Client:            | Ardmore Chemical  |                     | Date Collected: |               |
| Project:           | PVSC Monthly 2025 |                     | Date Received:  |               |
| Client Sample ID:  | PB170210BL        |                     | SDG No.:        | Q3385         |
| Lab Sample ID:     | PB170210BL        |                     | Matrix:         | Water         |
| Analytical Method: | 625.1             | Level : LOW         | % Solid:        | 0             |
| Sample Wt/Vol:     | 1000 mL           | Final Vol: 1000 uL  | Test:           | SVOCMS Group1 |
| Prep Method :      | 3510C             | Prep Date: 10/22/25 |                 |               |

| CAS Number | Parameter                  | Conc. | Qua. | DF | MDL  | LOQ / CRQL | Units | Date Ana.      | Prep BatchID |
|------------|----------------------------|-------|------|----|------|------------|-------|----------------|--------------|
| 85-68-7    | Butylbenzylphthalate       | 1.90  | U    | 1  | 1.90 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 91-94-1    | 3,3-Dichlorobenzidine      | 0.93  | U    | 1  | 0.93 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 56-55-3    | Benzo(a)anthracene         | 0.45  | U    | 1  | 0.45 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 218-01-9   | Chrysene                   | 0.44  | U    | 1  | 0.44 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 1.60  | U    | 1  | 1.60 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 117-84-0   | Di-n-octyl phthalate       | 2.30  | U    | 1  | 2.30 | 10.0       | ug/L  | 10/27/25 12:26 | PB170210     |
| 205-99-2   | Benzo(b)fluoranthene       | 0.49  | U    | 1  | 0.49 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 207-08-9   | Benzo(k)fluoranthene       | 0.48  | U    | 1  | 0.48 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 50-32-8    | Benzo(a)pyrene             | 0.55  | U    | 1  | 0.55 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 0.59  | U    | 1  | 0.59 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 53-70-3    | Dibenzo(a,h)anthracene     | 0.67  | U    | 1  | 0.67 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |
| 191-24-2   | Benzo(g,h,i)perylene       | 0.69  | U    | 1  | 0.69 | 5.00       | ug/L  | 10/27/25 12:26 | PB170210     |

### SURROGATES

|            |                      |      |  |  |          |      |          |
|------------|----------------------|------|--|--|----------|------|----------|
| 367-12-4   | 2-Fluorophenol       | 82.3 |  |  | 60 - 140 | 82%  | SPK: 100 |
| 13127-88-3 | Phenol-d6            | 87.7 |  |  | 60 - 140 | 88%  | SPK: 100 |
| 4165-60-0  | Nitrobenzene-d5      | 84.1 |  |  | 60 - 140 | 84%  | SPK: 100 |
| 321-60-8   | 2-Fluorobiphenyl     | 84.3 |  |  | 60 - 140 | 84%  | SPK: 100 |
| 118-79-6   | 2,4,6-Tribromophenol | 79.9 |  |  | 60 - 140 | 80%  | SPK: 100 |
| 1718-51-0  | Terphenyl-d14        | 104  |  |  | 60 - 140 | 104% | SPK: 100 |

### INTERNAL STANDARDS

| INTERNAL STANDARDS |                        | Area Count |
|--------------------|------------------------|------------|
| 3855-82-1          | 1,4-Dichlorobenzene-d4 | 69800      |
| 1146-65-2          | Naphthalene-d8         | 269000     |
| 15067-26-2         | Acenaphthene-d10       | 148000     |
| 1517-22-2          | Phenanthrene-d10       | 294000     |
| 1719-03-5          | Chrysene-d12           | 173000     |
| 1520-96-3          | Perylene-d12           | 183000     |

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