

## Report of Analysis

|                    |                   |                  |                 |               |
|--------------------|-------------------|------------------|-----------------|---------------|
| Client:            | Ardmore Chemical  |                  | Date Collected: |               |
| Project:           | PVSC Monthly 2025 |                  | Date Received:  |               |
| Client Sample ID:  | PB168904BS        |                  | SDG No.:        | Q2585         |
| Lab Sample ID:     | PB168904BS        |                  | Matrix:         | Water         |
| Analytical Method: | 625.1             |                  | % Solid:        | 0             |
| Sample Wt/Vol:     | 1000              | 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 |
| BF143162.D        | 1         | 07/17/25 08:39 | 07/18/25 12:53 | PB168904      |

| CAS Number     | Parameter                   | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|-----------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                             |       |           |      |            |       |
| 62-75-9        | n-Nitrosodimethylamine      | 42.4  |           | 0.86 | 10.0       | ug/L  |
| 108-95-2       | Phenol                      | 42.4  |           | 0.91 | 5.00       | ug/L  |
| 111-44-4       | bis(2-Chloroethyl)ether     | 42.5  |           | 0.81 | 5.00       | ug/L  |
| 95-57-8        | 2-Chlorophenol              | 42.7  |           | 0.58 | 5.00       | ug/L  |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 42.7  |           | 1.30 | 5.00       | ug/L  |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 42.4  |           | 1.40 | 5.00       | ug/L  |
| 67-72-1        | Hexachloroethane            | 43.2  |           | 0.65 | 5.00       | ug/L  |
| 98-95-3        | Nitrobenzene                | 44.0  |           | 0.76 | 5.00       | ug/L  |
| 78-59-1        | Isophorone                  | 43.8  |           | 0.75 | 5.00       | ug/L  |
| 88-75-5        | 2-Nitrophenol               | 46.4  |           | 1.80 | 5.00       | ug/L  |
| 105-67-9       | 2,4-Dimethylphenol          | 44.1  |           | 1.90 | 5.00       | ug/L  |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 43.2  |           | 0.68 | 5.00       | ug/L  |
| 120-83-2       | 2,4-Dichlorophenol          | 43.6  |           | 0.52 | 5.00       | ug/L  |
| 120-82-1       | 1,2,4-Trichlorobenzene      | 42.7  |           | 0.54 | 5.00       | ug/L  |
| 91-20-3        | Naphthalene                 | 42.9  |           | 0.50 | 5.00       | ug/L  |
| 87-68-3        | Hexachlorobutadiene         | 42.9  |           | 0.54 | 5.00       | ug/L  |
| 59-50-7        | 4-Chloro-3-methylphenol     | 44.1  |           | 0.59 | 5.00       | ug/L  |
| 77-47-4        | Hexachlorocyclopentadiene   | 86.3  | E         | 3.60 | 10.0       | ug/L  |
| 88-06-2        | 2,4,6-Trichlorophenol       | 42.9  |           | 0.51 | 5.00       | ug/L  |
| 91-58-7        | 2-Chloronaphthalene         | 42.6  |           | 0.61 | 5.00       | ug/L  |
| 131-11-3       | Dimethylphthalate           | 43.9  |           | 0.61 | 5.00       | ug/L  |
| 208-96-8       | Acenaphthylene              | 43.2  |           | 0.75 | 5.00       | ug/L  |
| 606-20-2       | 2,6-Dinitrotoluene          | 47.2  |           | 0.92 | 5.00       | ug/L  |
| 83-32-9        | Acenaphthene                | 46.2  |           | 0.55 | 5.00       | ug/L  |
| 51-28-5        | 2,4-Dinitrophenol           | 90.5  | E         | 6.00 | 10.0       | ug/L  |
| 100-02-7       | 4-Nitrophenol               | 89.2  | E         | 2.40 | 10.0       | ug/L  |
| 121-14-2       | 2,4-Dinitrotoluene          | 48.9  |           | 1.20 | 5.00       | ug/L  |
| 84-66-2        | Diethylphthalate            | 44.7  |           | 0.69 | 5.00       | ug/L  |
| 7005-72-3      | 4-Chlorophenyl-phenylether  | 43.4  |           | 0.68 | 5.00       | ug/L  |

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| Project:           | PVSC Monthly 2025 |                  | Date Received:  |               |      |
| Client Sample ID:  | PB168904BS        |                  | SDG No.:        | Q2585         |      |
| Lab Sample ID:     | PB168904BS        |                  | Matrix:         | Water         |      |
| Analytical Method: | 625.1             |                  | % Solid:        | 0             |      |
| Sample Wt/Vol:     | 1000              | 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 |
| BF143162.D        | 1         | 07/17/25 08:39 | 07/18/25 12:53 | PB168904      |

| CAS Number | Parameter                  | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|----------------------------|-------|-----------|------|------------|-------|
| 86-73-7    | Fluorene                   | 42.8  |           | 0.63 | 5.00       | ug/L  |
| 534-52-1   | 4,6-Dinitro-2-methylphenol | 45.1  |           | 2.90 | 10.0       | ug/L  |
| 86-30-6    | n-Nitrosodiphenylamine     | 43.8  |           | 0.58 | 5.00       | ug/L  |
| 103-33-3   | Azobenzene                 | 44.1  |           | 0.81 | 5.00       | ug/L  |
| 101-55-3   | 4-Bromophenyl-phenylether  | 44.1  |           | 0.40 | 5.00       | ug/L  |
| 118-74-1   | Hexachlorobenzene          | 44.7  |           | 0.52 | 5.00       | ug/L  |
| 87-86-5    | Pentachlorophenol          | 91.0  | E         | 1.60 | 10.0       | ug/L  |
| 85-01-8    | Phenanthrene               | 43.8  |           | 0.50 | 5.00       | ug/L  |
| 120-12-7   | Anthracene                 | 43.8  |           | 0.61 | 5.00       | ug/L  |
| 84-74-2    | Di-n-butylphthalate        | 46.0  |           | 1.20 | 5.00       | ug/L  |
| 206-44-0   | Fluoranthene               | 44.2  |           | 0.82 | 5.00       | ug/L  |
| 92-87-5    | Benzidine                  | 41.4  |           | 4.30 | 10.0       | ug/L  |
| 129-00-0   | Pyrene                     | 44.7  |           | 0.50 | 5.00       | ug/L  |
| 85-68-7    | Butylbenzylphthalate       | 47.3  |           | 1.90 | 5.00       | ug/L  |
| 91-94-1    | 3,3-Dichlorobenzidine      | 25.1  |           | 0.93 | 10.0       | ug/L  |
| 56-55-3    | Benzo(a)anthracene         | 44.2  |           | 0.45 | 5.00       | ug/L  |
| 218-01-9   | Chrysene                   | 45.1  |           | 0.44 | 5.00       | ug/L  |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 47.8  |           | 1.60 | 5.00       | ug/L  |
| 117-84-0   | Di-n-octyl phthalate       | 46.5  |           | 2.30 | 10.0       | ug/L  |
| 205-99-2   | Benzo(b)fluoranthene       | 44.7  |           | 0.49 | 5.00       | ug/L  |
| 207-08-9   | Benzo(k)fluoranthene       | 46.5  |           | 0.48 | 5.00       | ug/L  |
| 50-32-8    | Benzo(a)pyrene             | 45.7  |           | 0.55 | 5.00       | ug/L  |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 44.8  |           | 0.59 | 5.00       | ug/L  |
| 53-70-3    | Dibenzo(a,h)anthracene     | 45.0  |           | 0.67 | 5.00       | ug/L  |
| 191-24-2   | Benzo(g,h,i)perylene       | 44.9  |           | 0.69 | 5.00       | ug/L  |

### SURROGATES

|            |                  |      |  |          |     |          |
|------------|------------------|------|--|----------|-----|----------|
| 367-12-4   | 2-Fluorophenol   | 79.6 |  | 60 - 140 | 80% | SPK: 100 |
| 13127-88-3 | Phenol-d6        | 80.9 |  | 60 - 140 | 81% | SPK: 100 |
| 4165-60-0  | Nitrobenzene-d5  | 80.0 |  | 60 - 140 | 80% | SPK: 100 |
| 321-60-8   | 2-Fluorobiphenyl | 74.5 |  | 60 - 140 | 74% | SPK: 100 |

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| Sample Wt/Vol:     | 1000              | Units: mL        | Final Vol:      | 1000 uL       |
| Soil Aliquot Vol:  |                   | uL               | Test:           | SVOCMS Group1 |
| Extraction Type :  |                   | Decanted : N     | Level :         | LOW           |
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| Prep Method :      | 3510C             |                  |                 |               |

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| BF143162.D        | 1         | 07/17/25 08:39 | 07/18/25 12:53 | PB168904      |

| CAS Number | Parameter            | Conc. | Qualifier | MDL      | LOQ / CRQL | Units    |
|------------|----------------------|-------|-----------|----------|------------|----------|
| 118-79-6   | 2,4,6-Tribromophenol | 83.6  |           | 60 - 140 | 84%        | SPK: 100 |
| 1718-51-0  | Terphenyl-d14        | 78.8  |           | 60 - 140 | 79%        | SPK: 100 |

### INTERNAL STANDARDS

|            |                        |        |        |
|------------|------------------------|--------|--------|
| 3855-82-1  | 1,4-Dichlorobenzene-d4 | 141000 | 6.963  |
| 1146-65-2  | Naphthalene-d8         | 541000 | 8.245  |
| 15067-26-2 | Acenaphthene-d10       | 282000 | 9.998  |
| 1517-22-2  | Phenanthrene-d10       | 448000 | 11.48  |
| 1719-03-5  | Chrysene-d12           | 250000 | 14.121 |
| 1520-96-3  | Perylene-d12           | 283000 | 15.621 |

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