

## Report of Analysis

Client: PSEG  
Project: PSEG Ridgefield OH ROW Bridge  
Client Sample ID: FRAC-TANK-M21012  
Lab Sample ID: Q3660-01  
Analytical Method: 8270E Level: LOW  
Sample Wt/Vol: 970 mL Final Vol: 1000 uL  
Prep Method : 3510C Prep Date: 11/18/25

Date Collected: 11/17/25  
Date Received: 11/17/25  
SDG No.: Q3660  
Matrix: Water  
% Solid: 0  
Test: SVOC-TCL BNA -20

| CAS Number     | Parameter                   | Conc. | Qua. | DF | MDL  | LOQ / CRQL | Units | Date Ana.      | Prep BatchID |
|----------------|-----------------------------|-------|------|----|------|------------|-------|----------------|--------------|
| <b>TARGETS</b> |                             |       |      |    |      |            |       |                |              |
| 100-52-7       | Benzaldehyde                | 10.3  | U    | 1  | 4.00 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 108-95-2       | Phenol                      | 5.20  | U    | 1  | 0.94 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 111-44-4       | bis(2-Chloroethyl)ether     | 5.20  | U    | 1  | 0.84 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 95-57-8        | 2-Chlorophenol              | 5.20  | U    | 1  | 0.60 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 95-48-7        | 2-Methylphenol              | 5.20  | U    | 1  | 1.20 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 108-60-1       | 2,2-oxybis(1-Chloropropane) | 5.20  | U    | 1  | 1.30 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 98-86-2        | Acetophenone                | 5.20  | U    | 1  | 0.76 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 65794-96-9     | 3+4-Methylphenols           | 10.3  | U    | 1  | 1.10 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 621-64-7       | n-Nitroso-di-n-propylamine  | 2.60  | U    | 1  | 1.50 | 2.60       | ug/L  | 11/18/25 18:23 | PB170608     |
| 67-72-1        | Hexachloroethane            | 5.20  | U    | 1  | 0.67 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 98-95-3        | Nitrobenzene                | 5.20  | U    | 1  | 0.78 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 78-59-1        | Isophorone                  | 5.20  | U    | 1  | 0.77 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 88-75-5        | 2-Nitrophenol               | 5.20  | U    | 1  | 1.80 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 105-67-9       | 2,4-Dimethylphenol          | 5.20  | U    | 1  | 1.90 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 111-91-1       | bis(2-Chloroethoxy)methane  | 5.20  | U    | 1  | 0.70 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 120-83-2       | 2,4-Dichlorophenol          | 5.20  | U    | 1  | 0.54 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 91-20-3        | Naphthalene                 | 5.20  | U    | 1  | 0.52 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 106-47-8       | 4-Chloroaniline             | 5.20  | U    | 1  | 0.87 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 87-68-3        | Hexachlorobutadiene         | 5.20  | U    | 1  | 0.56 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 105-60-2       | Caprolactam                 | 10.3  | U    | 1  | 1.20 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 59-50-7        | 4-Chloro-3-methylphenol     | 5.20  | U    | 1  | 0.61 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 91-57-6        | 2-Methylnaphthalene         | 5.20  | U    | 1  | 0.58 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 77-47-4        | Hexachlorocyclopentadiene   | 10.3  | U    | 1  | 3.70 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 88-06-2        | 2,4,6-Trichlorophenol       | 5.20  | U    | 1  | 0.53 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 95-95-4        | 2,4,5-Trichlorophenol       | 5.20  | U    | 1  | 0.64 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 92-52-4        | 1,1-Biphenyl                | 5.20  | U    | 1  | 0.55 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 91-58-7        | 2-Chloronaphthalene         | 5.20  | U    | 1  | 0.63 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 88-74-4        | 2-Nitroaniline              | 5.20  | U    | 1  | 1.30 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 131-11-3       | Dimethylphthalate           | 5.20  | U    | 1  | 0.63 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 208-96-8       | Acenaphthylene              | 5.20  | U    | 1  | 0.77 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 606-20-2       | 2,6-Dinitrotoluene          | 5.20  | U    | 1  | 0.95 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 99-09-2        | 3-Nitroaniline              | 5.20  | U    | 1  | 1.10 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 83-32-9        | Acenaphthene                | 5.20  | U    | 1  | 0.57 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 51-28-5        | 2,4-Dinitrophenol           | 10.3  | U    | 1  | 6.20 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 100-02-7       | 4-Nitrophenol               | 10.3  | U    | 1  | 2.50 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 132-64-9       | Dibenzofuran                | 5.20  | U    | 1  | 0.63 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 121-14-2       | 2,4-Dinitrotoluene          | 5.20  | U    | 1  | 1.30 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 84-66-2        | Diethylphthalate            | 5.20  | U    | 1  | 0.71 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 7005-72-3      | 4-Chlorophenyl-phenylether  | 5.20  | U    | 1  | 0.70 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 86-73-7        | Fluorene                    | 5.20  | U    | 1  | 0.65 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 100-01-6       | 4-Nitroaniline              | 5.20  | U    | 1  | 1.50 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 534-52-1       | 4,6-Dinitro-2-methylphenol  | 10.3  | U    | 1  | 3.00 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |

## Report of Analysis

|                    |                               |                 |                  |
|--------------------|-------------------------------|-----------------|------------------|
| Client:            | PSEG                          | Date Collected: | 11/17/25         |
| Project:           | PSEG Ridgefield OH ROW Bridge | Date Received:  | 11/17/25         |
| Client Sample ID:  | FRAC-TANK-M21012              | SDG No.:        | Q3660            |
| Lab Sample ID:     | Q3660-01                      | Matrix:         | Water            |
| Analytical Method: | 8270E                         | % Solid:        | 0                |
| Sample Wt/Vol:     | 970 mL                        | Test:           | SVOC-TCL BNA -20 |
| Prep Method :      | 3510C                         |                 |                  |
|                    | Level: LOW                    |                 |                  |
|                    | Final Vol: 1000 uL            |                 |                  |
|                    | Prep Date: 11/18/25           |                 |                  |

| CAS Number | Parameter                  | Conc. | Qua. | DF | MDL  | LOQ / CRQL | Units | Date Ana.      | Prep BatchID |
|------------|----------------------------|-------|------|----|------|------------|-------|----------------|--------------|
| 86-30-6    | n-Nitrosodiphenylamine     | 5.20  | U    | 1  | 0.60 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 101-55-3   | 4-Bromophenyl-phenylether  | 5.20  | U    | 1  | 0.41 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 118-74-1   | Hexachlorobenzene          | 5.20  | U    | 1  | 0.54 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 1912-24-9  | Atrazine                   | 5.20  | U    | 1  | 1.00 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 87-86-5    | Pentachlorophenol          | 10.3  | U    | 1  | 1.60 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 85-01-8    | Phenanthrene               | 5.20  | U    | 1  | 0.52 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 120-12-7   | Anthracene                 | 5.20  | U    | 1  | 0.63 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 86-74-8    | Carbazole                  | 5.20  | U    | 1  | 0.74 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 84-74-2    | Di-n-butylphthalate        | 5.20  | U    | 1  | 1.30 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 206-44-0   | Fluoranthene               | 5.20  | U    | 1  | 0.85 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 129-00-0   | Pyrene                     | 5.20  | U    | 1  | 0.52 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 85-68-7    | Butylbenzylphthalate       | 5.20  | U    | 1  | 2.00 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 91-94-1    | 3,3-Dichlorobenzidine      | 10.3  | U    | 1  | 0.96 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 56-55-3    | Benzo(a)anthracene         | 5.20  | U    | 1  | 0.46 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 218-01-9   | Chrysene                   | 5.20  | U    | 1  | 0.45 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 117-81-7   | Bis(2-ethylhexyl)phthalate | 5.20  | U    | 1  | 1.60 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 117-84-0   | Di-n-octyl phthalate       | 10.3  | U    | 1  | 2.40 | 10.3       | ug/L  | 11/18/25 18:23 | PB170608     |
| 205-99-2   | Benzo(b)fluoranthene       | 5.20  | U    | 1  | 0.51 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 207-08-9   | Benzo(k)fluoranthene       | 5.20  | U    | 1  | 0.49 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 50-32-8    | Benzo(a)pyrene             | 5.20  | U    | 1  | 0.57 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 193-39-5   | Indeno(1,2,3-cd)pyrene     | 5.20  | U    | 1  | 0.61 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 53-70-3    | Dibenzo(a,h)anthracene     | 5.20  | U    | 1  | 0.69 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 191-24-2   | Benzo(g,h,i)perylene       | 5.20  | U    | 1  | 0.71 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 95-94-3    | 1,2,4,5-Tetrachlorobenzene | 5.20  | U    | 1  | 0.54 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 123-91-1   | 1,4-Dioxane                | 5.20  | U    | 1  | 1.00 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |
| 58-90-2    | 2,3,4,6-Tetrachlorophenol  | 5.20  | U    | 1  | 0.74 | 5.20       | ug/L  | 11/18/25 18:23 | PB170608     |

### SURROGATES

|            |                      |      |  |          |      |          |
|------------|----------------------|------|--|----------|------|----------|
| 367-12-4   | 2-Fluorophenol       | 66.9 |  | 10 - 134 | 45%  | SPK: 150 |
| 13127-88-3 | Phenol-d6            | 42.9 |  | 10 - 135 | 29%  | SPK: 150 |
| 4165-60-0  | Nitrobenzene-d5      | 86.6 |  | 39 - 138 | 87%  | SPK: 100 |
| 321-60-8   | 2-Fluorobiphenyl     | 86.2 |  | 52 - 132 | 86%  | SPK: 100 |
| 118-79-6   | 2,4,6-Tribromophenol | 124  |  | 44 - 137 | 83%  | SPK: 150 |
| 1718-51-0  | Terphenyl-d14        | 101  |  | 42 - 152 | 101% | SPK: 100 |

### INTERNAL STANDARDS

|            |                        |        |
|------------|------------------------|--------|
| 3855-82-1  | 1,4-Dichlorobenzene-d4 | 45100  |
| 1146-65-2  | Naphthalene-d8         | 178000 |
| 15067-26-2 | Acenaphthene-d10       | 148000 |
| 1517-22-2  | Phenanthrene-d10       | 311000 |
| 1719-03-5  | Chrysene-d12           | 300000 |
| 1520-96-3  | Perylene-d12           | 340000 |

### TENTATIVE IDENTIFIED COMPOUNDS

|          |                |      |   |      |      |
|----------|----------------|------|---|------|------|
| 100-51-6 | Benzyl Alcohol | 9.70 | J | 8.39 | ug/L |
| 65-85-0  | Benzoic acid   | 15.1 | J | 10.2 | ug/L |

## Report of Analysis

|                    |                               |                 |                  |
|--------------------|-------------------------------|-----------------|------------------|
| Client:            | PSEG                          | Date Collected: | 11/17/25         |
| Project:           | PSEG Ridgefield OH ROW Bridge | Date Received:  | 11/17/25         |
| Client Sample ID:  | FRAC-TANK-M21012              | SDG No.:        | Q3660            |
| Lab Sample ID:     | Q3660-01                      | Matrix:         | Water            |
| Analytical Method: | 8270E                         | % Solid:        | 0                |
|                    | Level: LOW                    | Test:           | SVOC-TCL BNA -20 |
| Sample Wt/Vol:     | 970 mL                        |                 |                  |
|                    | Final Vol: 1000 uL            |                 |                  |
| Prep Method :      | 3510C                         |                 |                  |
|                    | Prep Date: 11/18/25           |                 |                  |

| CAS Number  | Parameter                          | Conc. | Qua. | DF | MDL | LOQ / CRQL | Units | Date Ana. | Prep BatchID |
|-------------|------------------------------------|-------|------|----|-----|------------|-------|-----------|--------------|
| 002782-91-4 | Thiourea, tetramethyl-             | 2.90  | J    |    |     | 11.3       | ug/L  |           |              |
| 000095-16-9 | Benzothiazole                      | 2.50  | J    |    |     | 11.6       | ug/L  |           |              |
| 000621-59-0 | Benzaldehyde, 3-hydroxy-4-methoxy- | 3.70  | J    |    |     | 13.7       | ug/L  |           |              |
| 000101-83-7 | Cyclohexanamine, N-cyclohexyl-     | 10.1  | J    |    |     | 14.3       | ug/L  |           |              |

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