

## Report of Analysis

|                    |                                        |                  |                 |               |
|--------------------|----------------------------------------|------------------|-----------------|---------------|
| Client:            | Alliance Technical Group, LLC - Newark |                  | Date Collected: |               |
| Project:           | NJ Waste Water PT                      |                  | Date Received:  |               |
| Client Sample ID:  | PB168238BSD                            |                  | SDG No.:        | Q2181         |
| Lab Sample ID:     | PB168238BSD                            |                  | Matrix:         | Water         |
| Analytical Method: | SW8270ESIM                             |                  | % Solid:        | 0             |
| Sample Wt/Vol:     | 1000                                   | Units: mL        | Final Vol:      | 1000 uL       |
| Soil Aliquot Vol:  |                                        | uL               | Test:           | SVOCMS Group5 |
| Extraction Type :  |                                        | Decanted : N     | Level :         | LOW           |
| Injection Volume : |                                        | GPC Factor : 1.0 | GPC Cleanup :   | N PH :        |
| Prep Method :      |                                        |                  |                 |               |

|                   |           |                |                |               |
|-------------------|-----------|----------------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date      | Date Analyzed  | Prep Batch ID |
| BN037169.D        | 1         | 06/02/25 08:55 | 06/04/25 13:03 | PB168238      |

| CAS Number                | Parameter               | Conc. | Qualifier | MDL      | LOQ / CRQL | Units    |
|---------------------------|-------------------------|-------|-----------|----------|------------|----------|
| <b>TARGETS</b>            |                         |       |           |          |            |          |
| 208-96-8                  | Acenaphthylene          | 0.39  |           | 0.040    | 0.10       | ug/L     |
| <b>SURROGATES</b>         |                         |       |           |          |            |          |
| 7297-45-2                 | 2-Methylnaphthalene-d10 | 0.38  |           | 20 - 139 | 96%        | SPK: 0.4 |
| 93951-69-0                | Fluoranthene-d10        | 0.30  |           | 54 - 157 | 74%        | SPK: 0.4 |
| 4165-60-0                 | Nitrobenzene-d5         | 0.36  |           | 27 - 154 | 90%        | SPK: 0.4 |
| 321-60-8                  | 2-Fluorobiphenyl        | 0.37  |           | 30 - 155 | 91%        | SPK: 0.4 |
| 1718-51-0                 | Terphenyl-d14           | 0.36  |           | 54 - 175 | 90%        | SPK: 0.4 |
| <b>INTERNAL STANDARDS</b> |                         |       |           |          |            |          |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4  | 2150  | 7.59      |          |            |          |
| 1146-65-2                 | Naphthalene-d8          | 5350  | 10.372    |          |            |          |
| 15067-26-2                | Acenaphthene-d10        | 2660  | 14.235    |          |            |          |
| 1517-22-2                 | Phenanthrene-d10        | 4420  | 16.984    |          |            |          |
| 1719-03-5                 | Chrysene-d12            | 2670  | 21.189    |          |            |          |
| 1520-96-3                 | Perylene-d12            | 2550  | 23.377    |          |            |          |

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