

## Report of Analysis

|                    |                                   |                 |                      |
|--------------------|-----------------------------------|-----------------|----------------------|
| Client:            | PSEG                              | Date Collected: | 03/17/25             |
| Project:           | Burlington Gas District           | Date Received:  | 03/17/25             |
| Client Sample ID:  | OILY-DEBRIS-COMP                  | SDG No.:        | Q1592                |
| Lab Sample ID:     | Q1592-01                          | Matrix:         | SOIL                 |
| Analytical Method: | SW8260                            | % Solid:        | 85.6                 |
| Sample Wt/Vol:     | 5.01      Units:    g             | Final Vol:      | 5000              uL |
| Soil Aliquot Vol:  | uL                                | Test:           | VOC-TCLVOA-10        |
| GC Column:         | RXI-624              ID :    0.25 | Level :         | LOW                  |
| Prep Method :      |                                   |                 |                      |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VY021565.D        | 1         |           | 03/19/25 12:40 | VY031925      |

| CAS Number     | Parameter                      | Conc.  | Qualifier | MDL     | LOQ / CRQL | Units(Dry Weight) |
|----------------|--------------------------------|--------|-----------|---------|------------|-------------------|
| <b>TARGETS</b> |                                |        |           |         |            |                   |
| 75-71-8        | Dichlorodifluoromethane        | 0.0058 | U         | 0.0013  | 0.0058     | mg/Kg             |
| 74-87-3        | Chloromethane                  | 0.0058 | U         | 0.0013  | 0.0058     | mg/Kg             |
| 75-01-4        | Vinyl Chloride                 | 0.0058 | U         | 0.00092 | 0.0058     | mg/Kg             |
| 74-83-9        | Bromomethane                   | 0.0058 | U         | 0.0012  | 0.0058     | mg/Kg             |
| 75-00-3        | Chloroethane                   | 0.0058 | U         | 0.0015  | 0.0058     | mg/Kg             |
| 75-69-4        | Trichlorofluoromethane         | 0.0058 | U         | 0.0014  | 0.0058     | mg/Kg             |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.0058 | U         | 0.0012  | 0.0058     | mg/Kg             |
| 75-35-4        | 1,1-Dichloroethene             | 0.0058 | U         | 0.0012  | 0.0058     | mg/Kg             |
| 67-64-1        | Acetone                        | 0.029  | U         | 0.0055  | 0.029      | mg/Kg             |
| 75-15-0        | Carbon Disulfide               | 0.0058 | U         | 0.0012  | 0.0058     | mg/Kg             |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.0058 | U         | 0.00085 | 0.0058     | mg/Kg             |
| 79-20-9        | Methyl Acetate                 | 0.0058 | U         | 0.0018  | 0.0058     | mg/Kg             |
| 75-09-2        | Methylene Chloride             | 0.012  | U         | 0.0041  | 0.012      | mg/Kg             |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.0058 | U         | 0.0010  | 0.0058     | mg/Kg             |
| 75-34-3        | 1,1-Dichloroethane             | 0.0058 | U         | 0.00093 | 0.0058     | mg/Kg             |
| 110-82-7       | Cyclohexane                    | 0.0058 | U         | 0.00092 | 0.0058     | mg/Kg             |
| 78-93-3        | 2-Butanone                     | 0.029  | U         | 0.0076  | 0.029      | mg/Kg             |
| 56-23-5        | Carbon Tetrachloride           | 0.0058 | U         | 0.0011  | 0.0058     | mg/Kg             |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.0058 | U         | 0.00087 | 0.0058     | mg/Kg             |
| 74-97-5        | Bromochloromethane             | 0.0058 | U         | 0.0013  | 0.0058     | mg/Kg             |
| 67-66-3        | Chloroform                     | 0.0058 | U         | 0.00098 | 0.0058     | mg/Kg             |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.0058 | U         | 0.0011  | 0.0058     | mg/Kg             |
| 108-87-2       | Methylcyclohexane              | 0.0058 | U         | 0.0011  | 0.0058     | mg/Kg             |
| 71-43-2        | Benzene                        | 0.0058 | U         | 0.00092 | 0.0058     | mg/Kg             |
| 107-06-2       | 1,2-Dichloroethane             | 0.0058 | U         | 0.00092 | 0.0058     | mg/Kg             |
| 79-01-6        | Trichloroethene                | 0.0058 | U         | 0.00094 | 0.0058     | mg/Kg             |
| 78-87-5        | 1,2-Dichloropropane            | 0.0058 | U         | 0.0011  | 0.0058     | mg/Kg             |
| 75-27-4        | Bromodichloromethane           | 0.0058 | U         | 0.00091 | 0.0058     | mg/Kg             |
| 108-10-1       | 4-Methyl-2-Pentanone           | 0.029  | U         | 0.0042  | 0.029      | mg/Kg             |
| 108-88-3       | Toluene                        | 0.0058 | U         | 0.00091 | 0.0058     | mg/Kg             |

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| Soil Aliquot Vol:  | uL                                | Test:           | VOC-TCLVOA-10        |
| GC Column:         | RXI-624              ID :    0.25 | Level :         | LOW                  |
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| VY021565.D        | 1         |           | 03/19/25 12:40 | VY031925      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units(Dry Weight) |
|---------------------------|-----------------------------|--------|-----------|----------|------------|-------------------|
| 10061-02-6                | t-1,3-Dichloropropene       | 0.0058 | U         | 0.00076  | 0.0058     | mg/Kg             |
| 10061-01-5                | cis-1,3-Dichloropropene     | 0.0058 | U         | 0.00072  | 0.0058     | mg/Kg             |
| 79-00-5                   | 1,1,2-Trichloroethane       | 0.0058 | U         | 0.0011   | 0.0058     | mg/Kg             |
| 591-78-6                  | 2-Hexanone                  | 0.029  | U         | 0.0043   | 0.029      | mg/Kg             |
| 124-48-1                  | Dibromochloromethane        | 0.0058 | U         | 0.0010   | 0.0058     | mg/Kg             |
| 106-93-4                  | 1,2-Dibromoethane           | 0.0058 | U         | 0.0010   | 0.0058     | mg/Kg             |
| 127-18-4                  | Tetrachloroethene           | 0.0058 | U         | 0.0012   | 0.0058     | mg/Kg             |
| 108-90-7                  | Chlorobenzene               | 0.0058 | U         | 0.0011   | 0.0058     | mg/Kg             |
| 100-41-4                  | Ethyl Benzene               | 0.0058 | U         | 0.00078  | 0.0058     | mg/Kg             |
| 179601-23-1               | m/p-Xylenes                 | 0.012  | U         | 0.0014   | 0.012      | mg/Kg             |
| 95-47-6                   | o-Xylene                    | 0.0058 | U         | 0.00096  | 0.0058     | mg/Kg             |
| 100-42-5                  | Styrene                     | 0.0058 | U         | 0.00083  | 0.0058     | mg/Kg             |
| 75-25-2                   | Bromoform                   | 0.0058 | U         | 0.0010   | 0.0058     | mg/Kg             |
| 98-82-8                   | Isopropylbenzene            | 0.0058 | U         | 0.00091  | 0.0058     | mg/Kg             |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 0.0058 | U         | 0.0014   | 0.0058     | mg/Kg             |
| 541-73-1                  | 1,3-Dichlorobenzene         | 0.0058 | U         | 0.0020   | 0.0058     | mg/Kg             |
| 106-46-7                  | 1,4-Dichlorobenzene         | 0.0058 | U         | 0.0018   | 0.0058     | mg/Kg             |
| 95-50-1                   | 1,2-Dichlorobenzene         | 0.0058 | U         | 0.0017   | 0.0058     | mg/Kg             |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 0.0058 | U         | 0.0021   | 0.0058     | mg/Kg             |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 0.0058 | U         | 0.0035   | 0.0058     | mg/Kg             |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 0.0058 | U         | 0.0037   | 0.0058     | mg/Kg             |
| <b>SURROGATES</b>         |                             |        |           |          |            |                   |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 54.7   |           | 63 - 155 | 109%       | SPK: 50           |
| 1868-53-7                 | Dibromofluoromethane        | 51.3   |           | 70 - 134 | 103%       | SPK: 50           |
| 2037-26-5                 | Toluene-d8                  | 48.7   |           | 74 - 123 | 97%        | SPK: 50           |
| 460-00-4                  | 4-Bromofluorobenzene        | 41.3   |           | 38 - 136 | 83%        | SPK: 50           |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |                   |
| 363-72-4                  | Pentafluorobenzene          | 253000 | 7.707     |          |            |                   |
| 540-36-3                  | 1,4-Difluorobenzene         | 466000 | 8.615     |          |            |                   |
| 3114-55-4                 | Chlorobenzene-d5            | 407000 | 11.42     |          |            |                   |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 146000 | 13.352    |          |            |                   |

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|------------|-----------|-------|-----------|-----|------------|-------|
|------------|-----------|-------|-----------|-----|------------|-------|

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