

## Report of Analysis

|                    |                                   |                 |                      |
|--------------------|-----------------------------------|-----------------|----------------------|
| Client:            | PSEG                              | Date Collected: | 05/20/25             |
| Project:           | Waldwick Substation               | Date Received:  | 05/20/25             |
| Client Sample ID:  | RBR200044                         | SDG No.:        | Q2097                |
| Lab Sample ID:     | Q2097-03                          | Matrix:         | SOIL                 |
| Analytical Method: | 8260D                             | % Solid:        | 89.1                 |
| Sample Wt/Vol:     | 5.69      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 |
| VY022361.D        | 1         |           | 05/21/25 12:47 | VY052125      |

| CAS Number     | Parameter                      | Conc.  | Qualifier | MDL     | LOQ / CRQL | Units(Dry Weight) |
|----------------|--------------------------------|--------|-----------|---------|------------|-------------------|
| <b>TARGETS</b> |                                |        |           |         |            |                   |
| 75-71-8        | Dichlorodifluoromethane        | 0.0049 | U         | 0.0011  | 0.0049     | mg/Kg             |
| 74-87-3        | Chloromethane                  | 0.0049 | U         | 0.0011  | 0.0049     | mg/Kg             |
| 75-01-4        | Vinyl Chloride                 | 0.0049 | U         | 0.00078 | 0.0049     | mg/Kg             |
| 74-83-9        | Bromomethane                   | 0.0049 | U         | 0.0011  | 0.0049     | mg/Kg             |
| 75-00-3        | Chloroethane                   | 0.0049 | U         | 0.0012  | 0.0049     | mg/Kg             |
| 75-69-4        | Trichlorofluoromethane         | 0.0049 | U         | 0.0012  | 0.0049     | mg/Kg             |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.0049 | U         | 0.0010  | 0.0049     | mg/Kg             |
| 75-35-4        | 1,1-Dichloroethene             | 0.0049 | U         | 0.00099 | 0.0049     | mg/Kg             |
| 67-64-1        | Acetone                        | 0.025  | U         | 0.0047  | 0.025      | mg/Kg             |
| 75-15-0        | Carbon Disulfide               | 0.0049 | U         | 0.0010  | 0.0049     | mg/Kg             |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.0049 | U         | 0.00072 | 0.0049     | mg/Kg             |
| 79-20-9        | Methyl Acetate                 | 0.0049 | U         | 0.0015  | 0.0049     | mg/Kg             |
| 75-09-2        | Methylene Chloride             | 0.0099 | U         | 0.0035  | 0.0099     | mg/Kg             |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.0049 | U         | 0.00085 | 0.0049     | mg/Kg             |
| 75-34-3        | 1,1-Dichloroethane             | 0.0049 | U         | 0.00079 | 0.0049     | mg/Kg             |
| 110-82-7       | Cyclohexane                    | 0.0049 | U         | 0.00078 | 0.0049     | mg/Kg             |
| 78-93-3        | 2-Butanone                     | 0.025  | U         | 0.0064  | 0.025      | mg/Kg             |
| 56-23-5        | Carbon Tetrachloride           | 0.0049 | U         | 0.00096 | 0.0049     | mg/Kg             |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.0049 | U         | 0.00074 | 0.0049     | mg/Kg             |
| 74-97-5        | Bromochloromethane             | 0.0049 | U         | 0.0011  | 0.0049     | mg/Kg             |
| 67-66-3        | Chloroform                     | 0.0049 | U         | 0.00083 | 0.0049     | mg/Kg             |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.0049 | U         | 0.00092 | 0.0049     | mg/Kg             |
| 108-87-2       | Methylcyclohexane              | 0.0049 | U         | 0.00090 | 0.0049     | mg/Kg             |
| 71-43-2        | Benzene                        | 0.0049 | U         | 0.00078 | 0.0049     | mg/Kg             |
| 107-06-2       | 1,2-Dichloroethane             | 0.0049 | U         | 0.00078 | 0.0049     | mg/Kg             |
| 79-01-6        | Trichloroethene                | 0.0049 | U         | 0.00080 | 0.0049     | mg/Kg             |
| 78-87-5        | 1,2-Dichloropropane            | 0.0049 | U         | 0.00090 | 0.0049     | mg/Kg             |
| 75-27-4        | Bromodichloromethane           | 0.0049 | U         | 0.00077 | 0.0049     | mg/Kg             |
| 108-10-1       | 4-Methyl-2-Pentanone           | 0.025  | U         | 0.0035  | 0.025      | mg/Kg             |
| 108-88-3       | Toluene                        | 0.0049 | U         | 0.00077 | 0.0049     | mg/Kg             |

## Report of Analysis

|                    |                     |                 |               |
|--------------------|---------------------|-----------------|---------------|
| Client:            | PSEG                | Date Collected: | 05/20/25      |
| Project:           | Waldwick Substation | Date Received:  | 05/20/25      |
| Client Sample ID:  | RBR200044           | SDG No.:        | Q2097         |
| Lab Sample ID:     | Q2097-03            | Matrix:         | SOIL          |
| Analytical Method: | 8260D               | % Solid:        | 89.1          |
| Sample Wt/Vol:     | 5.69                | Units:          | g             |
| Soil Aliquot Vol:  |                     | Final Vol:      | 5000 uL       |
|                    |                     | Test:           | VOC-TCLVOA-10 |
| GC Column:         | RXI-624             | Level :         | LOW           |
| Prep Method :      | ID : 0.25           |                 |               |

|                   |           |           |                |               |
|-------------------|-----------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Prep Date | Date Analyzed  | Prep Batch ID |
| VY022361.D        | 1         |           | 05/21/25 12:47 | VY052125      |

| CAS Number  | Parameter                   | Conc.  | Qualifier | MDL     | LOQ / CRQL | Units(Dry Weight) |
|-------------|-----------------------------|--------|-----------|---------|------------|-------------------|
| 10061-02-6  | t-1,3-Dichloropropene       | 0.0049 | U         | 0.00064 | 0.0049     | mg/Kg             |
| 10061-01-5  | cis-1,3-Dichloropropene     | 0.0049 | U         | 0.00061 | 0.0049     | mg/Kg             |
| 79-00-5     | 1,1,2-Trichloroethane       | 0.0049 | U         | 0.00091 | 0.0049     | mg/Kg             |
| 591-78-6    | 2-Hexanone                  | 0.025  | U         | 0.0036  | 0.025      | mg/Kg             |
| 124-48-1    | Dibromochloromethane        | 0.0049 | U         | 0.00086 | 0.0049     | mg/Kg             |
| 106-93-4    | 1,2-Dibromoethane           | 0.0049 | U         | 0.00087 | 0.0049     | mg/Kg             |
| 127-18-4    | Tetrachloroethene           | 0.0049 | U         | 0.0010  | 0.0049     | mg/Kg             |
| 108-90-7    | Chlorobenzene               | 0.0049 | U         | 0.00090 | 0.0049     | mg/Kg             |
| 100-41-4    | Ethyl Benzene               | 0.0049 | U         | 0.00066 | 0.0049     | mg/Kg             |
| 179601-23-1 | m/p-Xylenes                 | 0.0099 | U         | 0.0012  | 0.0099     | mg/Kg             |
| 95-47-6     | o-Xylene                    | 0.0049 | U         | 0.00081 | 0.0049     | mg/Kg             |
| 100-42-5    | Styrene                     | 0.0049 | U         | 0.00070 | 0.0049     | mg/Kg             |
| 75-25-2     | Bromoform                   | 0.0049 | U         | 0.00085 | 0.0049     | mg/Kg             |
| 98-82-8     | Isopropylbenzene            | 0.0049 | U         | 0.00077 | 0.0049     | mg/Kg             |
| 79-34-5     | 1,1,2,2-Tetrachloroethane   | 0.0049 | U         | 0.0012  | 0.0049     | mg/Kg             |
| 541-73-1    | 1,3-Dichlorobenzene         | 0.0049 | U         | 0.0017  | 0.0049     | mg/Kg             |
| 106-46-7    | 1,4-Dichlorobenzene         | 0.0049 | U         | 0.0015  | 0.0049     | mg/Kg             |
| 95-50-1     | 1,2-Dichlorobenzene         | 0.0049 | U         | 0.0014  | 0.0049     | mg/Kg             |
| 96-12-8     | 1,2-Dibromo-3-Chloropropane | 0.0049 | U         | 0.0018  | 0.0049     | mg/Kg             |
| 120-82-1    | 1,2,4-Trichlorobenzene      | 0.0049 | U         | 0.0029  | 0.0049     | mg/Kg             |
| 87-61-6     | 1,2,3-Trichlorobenzene      | 0.0049 | U         | 0.0031  | 0.0049     | mg/Kg             |

## SURROGATES

|            |                       |      |          |     |         |
|------------|-----------------------|------|----------|-----|---------|
| 17060-07-0 | 1,2-Dichloroethane-d4 | 45.7 | 63 - 155 | 91% | SPK: 50 |
| 1868-53-7  | Dibromofluoromethane  | 49.2 | 70 - 134 | 98% | SPK: 50 |
| 2037-26-5  | Toluene-d8            | 49.0 | 74 - 123 | 98% | SPK: 50 |
| 460-00-4   | 4-Bromofluorobenzene  | 39.1 | 38 - 136 | 78% | SPK: 50 |

## INTERNAL STANDARDS

|           |                        |        |        |
|-----------|------------------------|--------|--------|
| 363-72-4  | Pentafluorobenzene     | 242000 | 7.719  |
| 540-36-3  | 1,4-Difluorobenzene    | 434000 | 8.622  |
| 3114-55-4 | Chlorobenzene-d5       | 346000 | 11.42  |
| 3855-82-1 | 1,4-Dichlorobenzene-d4 | 127000 | 13.352 |

### TENTATIVE IDENTIFIED COMPOUNDS

## Report of Analysis

|                    |                                           |                 |                              |
|--------------------|-------------------------------------------|-----------------|------------------------------|
| Client:            | PSEG                                      | Date Collected: | 05/20/25                     |
| Project:           | Waldwick Substation                       | Date Received:  | 05/20/25                     |
| Client Sample ID:  | RBR200044                                 | SDG No.:        | Q2097                        |
| Lab Sample ID:     | Q2097-03                                  | Matrix:         | SOIL                         |
| Analytical Method: | 8260D                                     | % Solid:        | 89.1                         |
| Sample Wt/Vol:     | 5.69      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 |
| VY022361.D        | 1         |           | 05/21/25 12:47 | VY052125      |

| CAS Number  | Parameter                          | Conc. | Qualifier | MDL | LOQ / CRQL | Units(Dry Weight) |
|-------------|------------------------------------|-------|-----------|-----|------------|-------------------|
| 054832-83-6 | 1H-Indene, octahydro-2,2,4,4,7,7-h | 14.8  | J         |     | 13.2       | ug/Kg             |

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