

## Report of Analysis

|                    |                                                    |           |                 |               |    |
|--------------------|----------------------------------------------------|-----------|-----------------|---------------|----|
| Client:            | Langan Engineering and Environmental Services, Inc |           | Date Collected: | 10/07/25      |    |
| Project:           | Con Edison Richmond Terrace                        |           | Date Received:  | 10/08/25      |    |
| Client Sample ID:  | SS-01-1-2                                          |           | SDG No.:        | Q3310         |    |
| Lab Sample ID:     | Q3310-02                                           |           | Matrix:         | SOIL          |    |
| Analytical Method: | 8260D                                              |           | % Solid:        | 72.4          |    |
| Sample Wt/Vol:     | 7.82                                               | 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: | Date Analyzed  | Prep Batch ID |
| VY023417.D        | 1         | 10/09/25 12:57 | VY100925      |

| CAS Number     | Parameter                      | Conc. | Qualifier | MDL  | LOQ / CRQL | Units(Dry Weight) |
|----------------|--------------------------------|-------|-----------|------|------------|-------------------|
| <b>TARGETS</b> |                                |       |           |      |            |                   |
| 75-71-8        | Dichlorodifluoromethane        | 1.00  | U         | 1.00 | 4.40       | ug/Kg             |
| 74-87-3        | Chloromethane                  | 1.00  | U         | 1.00 | 4.40       | ug/Kg             |
| 75-01-4        | Vinyl Chloride                 | 0.70  | U         | 0.70 | 4.40       | ug/Kg             |
| 74-83-9        | Bromomethane                   | 0.94  | U         | 0.94 | 4.40       | ug/Kg             |
| 75-00-3        | Chloroethane                   | 1.10  | U         | 1.10 | 4.40       | ug/Kg             |
| 75-69-4        | Trichlorofluoromethane         | 1.10  | U         | 1.10 | 4.40       | ug/Kg             |
| 76-13-1        | 1,1,2-Trichlorotrifluoroethane | 0.94  | U         | 0.94 | 4.40       | ug/Kg             |
| 75-35-4        | 1,1-Dichloroethene             | 0.88  | U         | 0.88 | 4.40       | ug/Kg             |
| 67-64-1        | Acetone                        | 4.20  | U         | 4.20 | 22.1       | ug/Kg             |
| 75-15-0        | Carbon Disulfide               | 0.94  | U         | 0.94 | 4.40       | ug/Kg             |
| 1634-04-4      | Methyl tert-butyl Ether        | 0.64  | U         | 0.64 | 4.40       | ug/Kg             |
| 79-20-9        | Methyl Acetate                 | 1.40  | U         | 1.40 | 4.40       | ug/Kg             |
| 75-09-2        | Methylene Chloride             | 4.50  | J         | 3.10 | 8.80       | ug/Kg             |
| 156-60-5       | trans-1,2-Dichloroethene       | 0.76  | U         | 0.76 | 4.40       | ug/Kg             |
| 75-34-3        | 1,1-Dichloroethane             | 0.71  | U         | 0.71 | 4.40       | ug/Kg             |
| 110-82-7       | Cyclohexane                    | 0.70  | U         | 0.70 | 4.40       | ug/Kg             |
| 78-93-3        | 2-Butanone                     | 5.80  | U         | 5.80 | 22.1       | ug/Kg             |
| 56-23-5        | Carbon Tetrachloride           | 0.86  | U         | 0.86 | 4.40       | ug/Kg             |
| 156-59-2       | cis-1,2-Dichloroethene         | 0.66  | U         | 0.66 | 4.40       | ug/Kg             |
| 74-97-5        | Bromochloromethane             | 1.00  | U         | 1.00 | 4.40       | ug/Kg             |
| 67-66-3        | Chloroform                     | 0.74  | U         | 0.74 | 4.40       | ug/Kg             |
| 71-55-6        | 1,1,1-Trichloroethane          | 0.82  | U         | 0.82 | 4.40       | ug/Kg             |
| 108-87-2       | Methylcyclohexane              | 0.80  | U         | 0.80 | 4.40       | ug/Kg             |
| 71-43-2        | Benzene                        | 0.70  | U         | 0.70 | 4.40       | ug/Kg             |
| 107-06-2       | 1,2-Dichloroethane             | 0.70  | U         | 0.70 | 4.40       | ug/Kg             |
| 79-01-6        | Trichloroethene                | 0.72  | U         | 0.72 | 4.40       | ug/Kg             |
| 78-87-5        | 1,2-Dichloropropane            | 0.80  | U         | 0.80 | 4.40       | ug/Kg             |
| 75-27-4        | Bromodichloromethane           | 0.69  | U         | 0.69 | 4.40       | ug/Kg             |
| 108-10-1       | 4-Methyl-2-Pentanone           | 3.20  | U         | 3.20 | 22.1       | ug/Kg             |
| 108-88-3       | Toluene                        | 0.69  | U         | 0.69 | 4.40       | ug/Kg             |

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| Soil Aliquot Vol:  |                                                    | uL        | Test:           | VOC-TCLVOA-10 |    |
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| VY023417.D        | 1         | 10/09/25 12:57 | VY100925      |

| CAS Number                | Parameter                   | Conc.   | Qualifier | MDL      | LOQ / CRQL | Units(Dry Weight) |
|---------------------------|-----------------------------|---------|-----------|----------|------------|-------------------|
| 10061-02-6                | t-1,3-Dichloropropene       | 0.57    | U         | 0.57     | 4.40       | ug/Kg             |
| 10061-01-5                | cis-1,3-Dichloropropene     | 0.55    | U         | 0.55     | 4.40       | ug/Kg             |
| 79-00-5                   | 1,1,2-Trichloroethane       | 0.81    | U         | 0.81     | 4.40       | ug/Kg             |
| 591-78-6                  | 2-Hexanone                  | 3.30    | U         | 3.30     | 22.1       | ug/Kg             |
| 124-48-1                  | Dibromochloromethane        | 0.77    | U         | 0.77     | 4.40       | ug/Kg             |
| 106-93-4                  | 1,2-Dibromoethane           | 0.78    | U         | 0.78     | 4.40       | ug/Kg             |
| 127-18-4                  | Tetrachloroethene           | 0.93    | U         | 0.93     | 4.40       | ug/Kg             |
| 108-90-7                  | Chlorobenzene               | 0.80    | U         | 0.80     | 4.40       | ug/Kg             |
| 100-41-4                  | Ethyl Benzene               | 0.59    | U         | 0.59     | 4.40       | ug/Kg             |
| 179601-23-1               | m/p-Xylenes                 | 2.10    | J         | 1.10     | 8.80       | ug/Kg             |
| 95-47-6                   | o-Xylene                    | 1.40    | J         | 0.72     | 4.40       | ug/Kg             |
| 100-42-5                  | Styrene                     | 0.63    | U         | 0.63     | 4.40       | ug/Kg             |
| 75-25-2                   | Bromoform                   | 0.76    | U         | 0.76     | 4.40       | ug/Kg             |
| 98-82-8                   | Isopropylbenzene            | 0.69    | U         | 0.69     | 4.40       | ug/Kg             |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.10    | U         | 1.10     | 4.40       | ug/Kg             |
| 541-73-1                  | 1,3-Dichlorobenzene         | 1.50    | U         | 1.50     | 4.40       | ug/Kg             |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.40    | U         | 1.40     | 4.40       | ug/Kg             |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.30    | U         | 1.30     | 4.40       | ug/Kg             |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.60    | U         | 1.60     | 4.40       | ug/Kg             |
| 120-82-1                  | 1,2,4-Trichlorobenzene      | 2.60    | U         | 2.60     | 4.40       | ug/Kg             |
| 87-61-6                   | 1,2,3-Trichlorobenzene      | 2.80    | U         | 2.80     | 4.40       | ug/Kg             |
| <b>SURROGATES</b>         |                             |         |           |          |            |                   |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 60.1    |           | 63 - 155 | 120%       | SPK: 50           |
| 1868-53-7                 | Dibromofluoromethane        | 53.4    |           | 70 - 134 | 107%       | SPK: 50           |
| 2037-26-5                 | Toluene-d8                  | 51.2    |           | 74 - 123 | 102%       | SPK: 50           |
| 460-00-4                  | 4-Bromofluorobenzene        | 40.5    |           | 17 - 146 | 81%        | SPK: 50           |
| <b>INTERNAL STANDARDS</b> |                             |         |           |          |            |                   |
| 363-72-4                  | Pentafluorobenzene          | 558000  | 7.713     |          |            |                   |
| 540-36-3                  | 1,4-Difluorobenzene         | 1090000 | 8.616     |          |            |                   |
| 3114-55-4                 | Chlorobenzene-d5            | 941000  | 11.42     |          |            |                   |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 284000  | 13.346    |          |            |                   |

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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