

## LAB CHRONICLE

|                 |                                    |                   |                       |
|-----------------|------------------------------------|-------------------|-----------------------|
| <b>OrderID:</b> | Q3254                              | <b>OrderDate:</b> | 9/30/2025 12:37:00 PM |
| <b>Client:</b>  | Lockwood, Kessler & Bartlett, Inc. | <b>Project:</b>   | Ansonia Landfill 2025 |
| <b>Contact:</b> | John Gerlach                       | <b>Location:</b>  | D31,VOA Ref. #3 Water |

| LabID    | ClientID   | Matrix | Test         | Method   | Sample Date | Prep Date | Anal Date | Received |
|----------|------------|--------|--------------|----------|-------------|-----------|-----------|----------|
| Q3254-01 | MW-1       | Water  | VOCMS Group1 | 8260-Low | 09/29/25    |           | 10/03/25  | 09/30/25 |
| Q3254-03 | MW-2       | Water  | VOCMS Group1 | 8260-Low | 09/29/25    |           | 10/03/25  | 09/30/25 |
| Q3254-05 | MW-3       | Water  | VOCMS Group1 | 8260-Low | 09/29/25    |           | 10/03/25  | 09/30/25 |
| Q3254-07 | MW-4       | Water  | VOCMS Group1 | 8260-Low | 09/29/25    |           | 10/03/25  | 09/30/25 |
| Q3254-09 | TRIP-BLANK | Water  | VOCMS Group1 | 8260-Low | 09/29/25    |           | 10/03/25  | 09/30/25 |

**Hit Summary Sheet**  
**SW-846**

**SDG No.:** Q3254

**Client:** Lockwood, Kessler & Bartlett, Inc.

| Sample ID                     | Client ID           | Matrix | Parameter                   | Concentration | C | MDL  | RDL  | Units |
|-------------------------------|---------------------|--------|-----------------------------|---------------|---|------|------|-------|
| <b>Client ID:</b><br>Q3254-03 | <b>MW-2</b><br>MW-2 | Water  | Chlorobenzene               | 0.97          | J | 0.12 | 1.00 | ug/L  |
|                               |                     |        | <b>Total Voc :</b>          | 0.97          |   |      |      |       |
|                               |                     |        | <b>Total Concentration:</b> | 0.97          |   |      |      |       |
| <b>Client ID:</b><br>Q3254-07 | <b>MW-4</b><br>MW-4 | Water  | Chlorobenzene               | 0.71          | J | 0.12 | 1.00 | ug/L  |
|                               |                     |        | <b>Total Voc :</b>          | 0.71          |   |      |      |       |
|                               |                     |        | <b>Total Concentration:</b> | 0.71          |   |      |      |       |



# SAMPLE DATA

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-1                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-01                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5      Units:    mL                | Final Vol:      | 5000      uL |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI      ID :    0.18         | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039252.D        | 1         | 10/03/25 16:11 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 1.00  | U         | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-1                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-01                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039252.D        | 1         | 10/03/25 16:11 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | UQ        | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 59.5   |           | 74 - 125 | 119%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 58.8   |           | 75 - 124 | 118%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.8   |           | 86 - 113 | 92%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.1   |           | 77 - 121 | 90%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 362000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 619000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 641000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 308000 | 11.175    |          |            |         |

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

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-2                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-03                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039253.D        | 1         | 10/03/25 16:34 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 0.97  | J         | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-2                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-03                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039253.D        | 1         | 10/03/25 16:34 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | UQ        | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 58.7   |           | 74 - 125 | 117%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 61.4   |           | 75 - 124 | 123%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 46.7   |           | 86 - 113 | 93%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 46.2   |           | 77 - 121 | 92%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 309000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 515000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 545000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 261000 | 11.172    |          |            |         |

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

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-3                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-05                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5      Units:    mL                | Final Vol:      | 5000      uL |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI      ID :    0.18         | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039254.D        | 1         | 10/03/25 16:56 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 1.00  | U         | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-3                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-05                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039254.D        | 1         | 10/03/25 16:56 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | UQ        | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 58.4   |           | 74 - 125 | 117%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 56.4   |           | 75 - 124 | 113%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.2   |           | 86 - 113 | 90%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.3   |           | 77 - 121 | 91%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 448000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 801000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 824000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 400000 | 11.172    |          |            |         |

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

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-4                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-07                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5      Units:    mL                | Final Vol:      | 5000      uL |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI      ID :    0.18         | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039255.D        | 1         | 10/03/25 17:19 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 0.71  | J         | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | MW-4                               | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-07                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039255.D        | 1         | 10/03/25 17:19 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | UQ        | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 57.5   |           | 74 - 125 | 115%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 57.4   |           | 75 - 124 | 115%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.6   |           | 86 - 113 | 91%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.1   |           | 77 - 121 | 90%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 393000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 674000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 694000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 336000 | 11.172    |          |            |         |

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

## Report of Analysis

|                    |                                    |                 |          |
|--------------------|------------------------------------|-----------------|----------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25 |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25 |
| Client Sample ID:  | TRIP-BLANK                         | SDG No.:        | Q3254    |
| Lab Sample ID:     | Q3254-09                           | Matrix:         | Water    |
| Analytical Method: | 8260D                              | % Solid:        | 0        |
| Sample Wt/Vol:     | 5                                  | Units:          | mL       |
| Soil Aliquot Vol:  |                                    |                 | uL       |
| GC Column:         | DB-624UI                           | ID :            | 0.18     |
| Prep Method :      |                                    | Level :         | LOW      |
|                    |                                    |                 |          |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039251.D        | 1         | 10/03/25 15:49 | VV100325      |

| CAS Number | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|------------|---------------------------|-------|-----------|------|------------|-------|
| TARGETS    |                           |       |           |      |            |       |
| 75-01-4    | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3    | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4    | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4    | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1   | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1    | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0    | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2    | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4   | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3    | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5    | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2   | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5    | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3    | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6    | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2    | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5    | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3    | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4    | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1   | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3   | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6 | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5 | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6   | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1   | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4   | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4   | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7   | Chlorobenzene             | 1.00  | U         | 0.12 | 1.00       | ug/L  |
| 630-20-6   | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4   | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |                 |              |
|--------------------|------------------------------------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. | Date Collected: | 09/29/25     |
| Project:           | Ansonia Landfill 2025              | Date Received:  | 09/30/25     |
| Client Sample ID:  | TRIP-BLANK                         | SDG No.:        | Q3254        |
| Lab Sample ID:     | Q3254-09                           | Matrix:         | Water        |
| Analytical Method: | 8260D                              | % Solid:        | 0            |
| Sample Wt/Vol:     | 5 Units: mL                        | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  | uL                                 | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI ID : 0.18                 | Level :         | LOW          |
| Prep Method :      |                                    |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039251.D        | 1         | 10/03/25 15:49 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | UQ        | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 58.2   |           | 74 - 125 | 116%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 58.9   |           | 75 - 124 | 118%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.8   |           | 86 - 113 | 92%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 44.6   |           | 77 - 121 | 89%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 330000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 558000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 580000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 279000 | 11.175    |          |            |         |

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



# QC SUMMARY

## Surrogate Summary

SDG No.: **Q3254**

Client: **Lockwood, Kessler & Bartlett, Inc.**

Analytical Method: **SW8260-Low**

| Lab Sample ID | Client ID    | Parameter             | Spike | Result | Recovery (%) | Qual | Limits (%) |      |
|---------------|--------------|-----------------------|-------|--------|--------------|------|------------|------|
|               |              |                       |       |        |              |      | Low        | High |
| Q3254-01      | MW-1         | 1,2-Dichloroethane-d4 | 50    | 59.5   | 119          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 58.8   | 118          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.8   | 92           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 45.0   | 90           |      | 77         | 121  |
| Q3254-03      | MW-2         | 1,2-Dichloroethane-d4 | 50    | 58.7   | 117          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 61.4   | 123          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 46.7   | 93           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 46.2   | 92           |      | 77         | 121  |
| Q3254-05      | MW-3         | 1,2-Dichloroethane-d4 | 50    | 58.4   | 117          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 56.4   | 113          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.2   | 90           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 45.3   | 91           |      | 77         | 121  |
| Q3254-07      | MW-4         | 1,2-Dichloroethane-d4 | 50    | 57.5   | 115          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 57.4   | 115          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.6   | 91           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 45.1   | 90           |      | 77         | 121  |
| Q3254-09      | TRIP-BLANK   | 1,2-Dichloroethane-d4 | 50    | 58.2   | 116          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 58.9   | 118          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.8   | 92           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 44.6   | 89           |      | 77         | 121  |
| VV1003WBL01   | VV1003WBL01  | 1,2-Dichloroethane-d4 | 50    | 51.7   | 103          |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 52.5   | 105          |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 44.6   | 89           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 45.2   | 90           |      | 77         | 121  |
| VV1003WBS01   | VV1003WBS01  | 1,2-Dichloroethane-d4 | 50    | 41.7   | 83           |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 49.6   | 99           |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.8   | 92           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 53.4   | 107          |      | 77         | 121  |
| VV1003WBSD0   | VV1003WBSD01 | 1,2-Dichloroethane-d4 | 50    | 42.8   | 86           |      | 74         | 125  |
|               |              | Dibromofluoromethane  | 50    | 48.6   | 97           |      | 75         | 124  |
|               |              | Toluene-d8            | 50    | 45.3   | 91           |      | 86         | 113  |
|               |              | 4-Bromofluorobenzene  | 50    | 53.2   | 106          |      | 77         | 121  |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

|                 |                                               |                           |                   |
|-----------------|-----------------------------------------------|---------------------------|-------------------|
| <b>SDG No.:</b> | <u>Q3254</u>                                  | <b>Analytical Method:</b> | <u>SW8260-Low</u> |
| <b>Client:</b>  | <u>Lockwood, Kessler &amp; Bartlett, Inc.</u> | <b>Datafile :</b>         | <u>VV039245.D</u> |

| Lab Sample ID | Parameter                   | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits<br>High | RPD |
|---------------|-----------------------------|-------|--------|------|-----|-----|------|-----|----------------|-----|
| VV1003WBS01   | Vinyl chloride              | 20    | 17.5   | ug/L | 88  |     |      | 65  | 117            |     |
|               | Chloroethane                | 20    | 18.1   | ug/L | 91  |     |      | 56  | 128            |     |
|               | Trichlorofluoromethane      | 20    | 19.1   | ug/L | 96  |     |      | 73  | 115            |     |
|               | 1,1-Dichloroethene          | 20    | 18.8   | ug/L | 94  |     |      | 74  | 110            |     |
|               | Acrylonitrile               | 100   | 88.9   | ug/L | 89  |     |      | 73  | 113            |     |
|               | Acetone                     | 100   | 91.0   | ug/L | 91  |     |      | 60  | 125            |     |
|               | Carbon disulfide            | 20    | 16.0   | ug/L | 80  |     |      | 64  | 112            |     |
|               | Methylene Chloride          | 20    | 19.9   | ug/L | 100 |     |      | 72  | 114            |     |
|               | Vinyl Acetate               | 100   | 88.0   | ug/L | 88  |     |      | 76  | 115            |     |
|               | 2-Butanone                  | 100   | 89.1   | ug/L | 89  |     |      | 65  | 122            |     |
|               | Carbon Tetrachloride        | 20    | 19.2   | ug/L | 96  |     |      | 77  | 113            |     |
|               | cis-1,2-Dichloroethene      | 20    | 19.2   | ug/L | 96  |     |      | 77  | 110            |     |
|               | Bromochloromethane          | 20    | 19.3   | ug/L | 97  |     |      | 70  | 124            |     |
|               | Chloroform                  | 20    | 18.1   | ug/L | 91  |     |      | 79  | 113            |     |
|               | 1,1,1-Trichloroethane       | 20    | 18.2   | ug/L | 91  |     |      | 80  | 108            |     |
|               | Benzene                     | 20    | 19.4   | ug/L | 97  |     |      | 82  | 109            |     |
|               | 1,2-Dichloropropane         | 20    | 19.9   | ug/L | 100 |     |      | 83  | 111            |     |
|               | Dibromomethane              | 20    | 19.8   | ug/L | 99  |     |      | 82  | 110            |     |
|               | Bromodichloromethane        | 20    | 19.3   | ug/L | 97  |     |      | 83  | 110            |     |
|               | 4-Methyl-2-Pentanone        | 100   | 100    | ug/L | 100 |     |      | 74  | 118            |     |
|               | Toluene                     | 20    | 20.9   | ug/L | 104 |     |      | 82  | 110            |     |
|               | t-1,3-Dichloropropene       | 20    | 21.1   | ug/L | 106 |     |      | 79  | 110            |     |
|               | cis-1,3-Dichloropropene     | 20    | 20.9   | ug/L | 104 |     |      | 82  | 110            |     |
|               | 2-Hexanone                  | 100   | 91.8   | ug/L | 92  |     |      | 73  | 117            |     |
|               | Dibromochloromethane        | 20    | 20.7   | ug/L | 104 |     |      | 82  | 110            |     |
|               | 1,2-Dibromoethane           | 20    | 20.6   | ug/L | 103 |     |      | 81  | 110            |     |
|               | Tetrachloroethene           | 20    | 20.3   | ug/L | 102 |     |      | 67  | 123            |     |
|               | Chlorobenzene               | 20    | 20.2   | ug/L | 101 |     |      | 82  | 109            |     |
|               | 1,1,1,2-Tetrachloroethane   | 20    | 20.3   | ug/L | 102 |     |      | 84  | 111            |     |
|               | Ethyl Benzene               | 20    | 20.2   | ug/L | 101 |     |      | 83  | 109            |     |
|               | m/p-Xylenes                 | 40    | 43.1   | ug/L | 108 |     |      | 82  | 110            |     |
|               | o-Xylene                    | 20    | 21.5   | ug/L | 108 |     |      | 83  | 109            |     |
|               | Styrene                     | 20    | 19.7   | ug/L | 99  |     |      | 80  | 111            |     |
|               | Bromoform                   | 20    | 21.5   | ug/L | 108 |     |      | 79  | 109            |     |
|               | 1,1,2,2-Tetrachloroethane   | 20    | 18.2   | ug/L | 91  |     |      | 76  | 118            |     |
|               | 1,2,3-Trichloropropane      | 20    | 18.1   | ug/L | 91  |     |      | 75  | 112            |     |
|               | 1,4-Dichlorobenzene         | 20    | 19.3   | ug/L | 97  |     |      | 82  | 107            |     |
|               | 1,2-Dichlorobenzene         | 20    | 19.7   | ug/L | 99  |     |      | 82  | 109            |     |
|               | 1,2-Dibromo-3-Chloropropane | 20    | 20.5   | ug/L | 103 |     |      | 68  | 112            |     |
|               | Methyl iodide               | 20    | 22.0   | ug/L | 110 |     |      | 70  | 130            |     |
|               | trans-1,4-Dichloro-2-butene | 20    | 19.0   | ug/L | 95  |     |      | 79  | 102            |     |

Laboratory Control Sample/Laboratory Control Sample Duplicate Summary

SW-846

|          |                                               |                    |                   |
|----------|-----------------------------------------------|--------------------|-------------------|
| SDG No.: | <u>Q3254</u>                                  | Analytical Method: | <u>SW8260-Low</u> |
| Client:  | <u>Lockwood, Kessler &amp; Bartlett, Inc.</u> | Datafile :         | <u>VV039246.D</u> |

| Lab Sample ID | Parameter                   | Spike | Result | Unit | Rec | RPD | Qual | Low | Limits<br>High | RPD |
|---------------|-----------------------------|-------|--------|------|-----|-----|------|-----|----------------|-----|
| VV1003WBSD01  | Vinyl chloride              | 20    | 17.0   | ug/L | 85  | 3   |      | 65  | 117            | 19  |
|               | Chloroethane                | 20    | 17.8   | ug/L | 89  | 2   |      | 56  | 128            | 20  |
|               | Trichlorofluoromethane      | 20    | 18.8   | ug/L | 94  | 2   |      | 73  | 115            | 16  |
|               | 1,1-Dichloroethene          | 20    | 18.7   | ug/L | 94  | 0   |      | 74  | 110            | 20  |
|               | Acrylonitrile               | 100   | 94.7   | ug/L | 95  | 7   |      | 73  | 113            | 29  |
|               | Acetone                     | 100   | 93.4   | ug/L | 93  | 2   |      | 60  | 125            | 20  |
|               | Carbon disulfide            | 20    | 15.9   | ug/L | 79  | 1   |      | 64  | 112            | 20  |
|               | Methylene Chloride          | 20    | 20.2   | ug/L | 101 | 1   |      | 72  | 114            | 20  |
|               | Vinyl Acetate               | 100   | 92.0   | ug/L | 92  | 4   |      | 76  | 115            | 26  |
|               | 2-Butanone                  | 100   | 93.5   | ug/L | 94  | 5   |      | 65  | 122            | 26  |
|               | Carbon Tetrachloride        | 20    | 18.8   | ug/L | 94  | 2   |      | 77  | 113            | 15  |
|               | cis-1,2-Dichloroethene      | 20    | 19.6   | ug/L | 98  | 2   |      | 77  | 110            | 20  |
|               | Bromochloromethane          | 20    | 19.6   | ug/L | 98  | 1   |      | 70  | 124            | 20  |
|               | Chloroform                  | 20    | 18.2   | ug/L | 91  | 0   |      | 79  | 113            | 20  |
|               | 1,1,1-Trichloroethane       | 20    | 18.4   | ug/L | 92  | 1   |      | 80  | 108            | 20  |
|               | Benzene                     | 20    | 19.3   | ug/L | 97  | 0   |      | 82  | 109            | 15  |
|               | 1,2-Dichloropropane         | 20    | 19.1   | ug/L | 96  | 4   |      | 83  | 111            | 16  |
|               | Dibromomethane              | 20    | 20.1   | ug/L | 101 | 2   |      | 82  | 110            | 20  |
|               | Bromodichloromethane        | 20    | 19.4   | ug/L | 97  | 0   |      | 83  | 110            | 16  |
|               | 4-Methyl-2-Pentanone        | 100   | 110    | ug/L | 110 | 10  |      | 74  | 118            | 25  |
|               | Toluene                     | 20    | 20.9   | ug/L | 104 | 0   |      | 82  | 110            | 16  |
|               | t-1,3-Dichloropropene       | 20    | 21.3   | ug/L | 106 | 0   |      | 79  | 110            | 20  |
|               | cis-1,3-Dichloropropene     | 20    | 21.1   | ug/L | 106 | 2   |      | 82  | 110            | 16  |
|               | 2-Hexanone                  | 100   | 95.2   | ug/L | 95  | 3   |      | 73  | 117            | 25  |
|               | Dibromochloromethane        | 20    | 20.6   | ug/L | 103 | 1   |      | 82  | 110            | 20  |
|               | 1,2-Dibromoethane           | 20    | 21.4   | ug/L | 107 | 4   |      | 81  | 110            | 20  |
|               | Tetrachloroethene           | 20    | 20.4   | ug/L | 102 | 0   |      | 67  | 123            | 15  |
|               | Chlorobenzene               | 20    | 20.2   | ug/L | 101 | 0   |      | 82  | 109            | 15  |
|               | 1,1,1,2-Tetrachloroethane   | 20    | 20.7   | ug/L | 104 | 2   |      | 84  | 111            | 20  |
|               | Ethyl Benzene               | 20    | 20.2   | ug/L | 101 | 0   |      | 83  | 109            | 16  |
|               | m/p-Xylenes                 | 40    | 43.3   | ug/L | 108 | 0   |      | 82  | 110            | 15  |
|               | o-Xylene                    | 20    | 21.8   | ug/L | 109 | 1   |      | 83  | 109            | 20  |
|               | Styrene                     | 20    | 19.6   | ug/L | 98  | 1   |      | 80  | 111            | 17  |
|               | Bromoform                   | 20    | 22.2   | ug/L | 111 | 3   | *    | 79  | 109            | 20  |
|               | 1,1,2,2-Tetrachloroethane   | 20    | 18.8   | ug/L | 94  | 3   |      | 76  | 118            | 20  |
|               | 1,2,3-Trichloropropane      | 20    | 18.8   | ug/L | 94  | 3   |      | 75  | 112            | 20  |
|               | 1,4-Dichlorobenzene         | 20    | 19.5   | ug/L | 98  | 1   |      | 82  | 107            | 15  |
|               | 1,2-Dichlorobenzene         | 20    | 20.2   | ug/L | 101 | 2   |      | 82  | 109            | 20  |
|               | 1,2-Dibromo-3-Chloropropane | 20    | 20.6   | ug/L | 103 | 0   |      | 68  | 112            | 20  |
|               | Methyl iodide               | 20    | 22.2   | ug/L | 111 | 1   |      | 70  | 130            | 20  |
|               | trans-1,4-Dichloro-2-butene | 20    | 19.5   | ug/L | 98  | 3   |      | 79  | 102            | 20  |

VOLATILE METHOD BLANK SUMMARY

Client ID

VV1003WBL01

Lab Name: Alliance

Contract: LOCK01

Lab Code: ACE

SDG NO.: Q3254

Lab File ID: VV039244.D

Lab Sample ID: VV1003WBL01

Date Analyzed: 10/03/2025

Time Analyzed: 11:26

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: (Y/N) N

Instrument ID: MSVOA\_V

THIS METHOD BLANK APPLIES TO THE FOLLOWING SAMPLES, MS AND MSD:

| EPA<br>SAMPLE NO. | LAB<br>SAMPLE ID | LAB<br>FILE ID | DATE<br>ANALYZED |
|-------------------|------------------|----------------|------------------|
| VV1003WBS01       | VV1003WBS01      | VV039245.D     | 10/03/2025       |
| VV1003WBSD01      | VV1003WBSD01     | VV039246.D     | 10/03/2025       |
| TRIP-BLANK        | Q3254-09         | VV039251.D     | 10/03/2025       |
| MW-1              | Q3254-01         | VV039252.D     | 10/03/2025       |
| MW-2              | Q3254-03         | VV039253.D     | 10/03/2025       |
| MW-3              | Q3254-05         | VV039254.D     | 10/03/2025       |
| MW-4              | Q3254-07         | VV039255.D     | 10/03/2025       |

COMMENTS: \_\_\_\_\_

\_\_\_\_\_

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: LOCK01

Lab Code: ACE

SDG NO.: Q3254

Lab File ID: VV039134.D

BFB Injection Date: 09/22/2025

Instrument ID: MSVOA\_V

BFB Injection Time: 10:15

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 19.6                 |
| 75  | 30.0 - 60.0% of mass 95            | 51.9                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 7.1                  |
| 173 | Less than 2.0% of mass 174         | 1.1 ( 1.6 ) 1        |
| 174 | 50.0 - 100.0% of mass 95           | 73.3                 |
| 175 | 5.0 - 9.0% of mass 174             | 5.5 ( 7.5 ) 1        |
| 176 | 95.0 - 101.0% of mass 174          | 70.4 ( 96 ) 1        |
| 177 | 5.0 - 9.0% of mass 176             | 4.4 ( 6.3 ) 2        |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| CLIENT ID   | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|-------------|---------------|-------------|---------------|---------------|
| VSTDICC010  | VSTDICC010    | VV039137.D  | 09/22/2025    | 12:26         |
| VSTDICC020  | VSTDICC020    | VV039138.D  | 09/22/2025    | 12:48         |
| VSTDICCC050 | VSTDICCC050   | VV039139.D  | 09/22/2025    | 13:26         |
| VSTDICC100  | VSTDICC100    | VV039140.D  | 09/22/2025    | 13:49         |
| VSTDICC005  | VSTDICC005    | VV039142.D  | 09/22/2025    | 15:37         |
| VSTDICC001  | VSTDICC001    | VV039143.D  | 09/22/2025    | 16:35         |

VOLATILE ORGANIC INSTRUMENT PERFORMANCE CHECK  
BROMOFLUOROBENZENE (BFB)

Lab Name: Alliance

Contract: LOCK01

Lab Code: ACE

SDG NO.: Q3254

Lab File ID: VV039241.D

BFB Injection Date: 10/03/2025

Instrument ID: MSVOA\_V

BFB Injection Time: 08:26

GC Column: DB-624UI ID: 0.18 (mm)

Heated Purge: Y/N N

| m/e | ION ABUNDANCE CRITERIA             | % RELATIVE ABUNDANCE |
|-----|------------------------------------|----------------------|
| 50  | 15.0 - 40.0% of mass 95            | 16.6                 |
| 75  | 30.0 - 60.0% of mass 95            | 49.5                 |
| 95  | Base Peak, 100% relative abundance | 100                  |
| 96  | 5.0 - 9.0% of mass 95              | 6.9                  |
| 173 | Less than 2.0% of mass 174         | 1 ( 1.3 ) 1          |
| 174 | 50.0 - 100.0% of mass 95           | 78                   |
| 175 | 5.0 - 9.0% of mass 174             | 5.5 ( 7 ) 1          |
| 176 | 95.0 - 101.0% of mass 174          | 74.9 ( 96 ) 1        |
| 177 | 5.0 - 9.0% of mass 176             | 5 ( 6.6 ) 2          |

1-Value is % mass 174

2-Value is % mass 176

THIS CHECK APPLIES TO THE FOLLOWING SAMPLES, MS, MSD, BLANKS, AND STANDARDS:

| CLIENT ID    | LAB SAMPLE ID | LAB FILE ID | DATE ANALYZED | TIME ANALYZED |
|--------------|---------------|-------------|---------------|---------------|
| VSTDCCC050   | VSTDCCC050    | VV039242.D  | 10/03/2025    | 08:53         |
| VV1003WBL01  | VV1003WBL01   | VV039244.D  | 10/03/2025    | 11:26         |
| VV1003WBS01  | VV1003WBS01   | VV039245.D  | 10/03/2025    | 13:07         |
| VV1003WBSD01 | VV1003WBSD01  | VV039246.D  | 10/03/2025    | 13:29         |
| TRIP-BLANK   | Q3254-09      | VV039251.D  | 10/03/2025    | 15:49         |
| MW-1         | Q3254-01      | VV039252.D  | 10/03/2025    | 16:11         |
| MW-2         | Q3254-03      | VV039253.D  | 10/03/2025    | 16:34         |
| MW-3         | Q3254-05      | VV039254.D  | 10/03/2025    | 16:56         |
| MW-4         | Q3254-07      | VV039255.D  | 10/03/2025    | 17:19         |

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

Lab Name: Alliance Contract: LOCK01  
 Lab Code: ACE SDG NO.: Q3254  
 Lab File ID: VV039242.D Date Analyzed: 10/03/2025  
 Instrument ID: MSVOA\_V Time Analyzed: 08:53  
 GC Column: DB-624UI ID: 0.18 (mm) Heated Purge: (Y/N) N

|                | IS1<br>AREA # | RT #  | IS2<br>AREA # | RT #  | IS3<br>AREA # | RT #  |
|----------------|---------------|-------|---------------|-------|---------------|-------|
| 12 HOUR STD    | 504801        | 4.60  | 717875        | 5.53  | 718198        | 8.77  |
| UPPER LIMIT    | 1009600       | 5.096 | 1435750       | 6.032 | 1436400       | 9.273 |
| LOWER LIMIT    | 252401        | 4.096 | 358938        | 5.032 | 359099        | 8.273 |
| EPA SAMPLE NO. |               |       |               |       |               |       |
| MW-1           | 361761        | 4.60  | 619136        | 5.54  | 640562        | 8.78  |
| MW-2           | 308662        | 4.60  | 515390        | 5.54  | 544820        | 8.78  |
| MW-3           | 448224        | 4.60  | 801027        | 5.54  | 823576        | 8.78  |
| MW-4           | 392891        | 4.60  | 674176        | 5.54  | 694020        | 8.78  |
| TRIP-BLANK     | 330074        | 4.60  | 558191        | 5.54  | 580452        | 8.78  |
| VV1003WBL01    | 501875        | 4.60  | 854024        | 5.53  | 854306        | 8.78  |
| VV1003WBS01    | 559480        | 4.60  | 866328        | 5.53  | 845304        | 8.77  |
| VV1003WBSD01   | 530712        | 4.60  | 834883        | 5.54  | 807929        | 8.78  |

IS1 = Pentafluorobenzene  
 IS2 = 1,4-Difluorobenzene  
 IS3 = Chlorobenzene-d5

AREA UPPER LIMIT = +100% of internal standard area  
 AREA LOWER LIMIT = -50% of internal standard area  
 RT UPPER LIMIT = +0.50 minutes of internal standard RT  
 RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.  
 \* Values outside of QC limits.

VOLATILE INTERNAL STANDARD AREA AND RT SUMMARY

|                |                   |                |                   |
|----------------|-------------------|----------------|-------------------|
| Lab Name:      | <u>Alliance</u>   | Contract:      | <u>LOCK01</u>     |
| Lab Code:      | <u>ACE</u>        | SDG NO.:       | <u>Q3254</u>      |
| Lab File ID:   | <u>VV039242.D</u> | Date Analyzed: | <u>10/03/2025</u> |
| Instrument ID: | <u>MSVOA_V</u>    | Time Analyzed: | <u>08:53</u>      |
| GC Column:     | <u>DB-624UI</u>   | ID:            | <u>0.18</u> (mm)  |
|                |                   | Heated Purge:  | (Y/N) <u>N</u>    |

|                | IS4<br>AREA # | RT #   |  |  |  |  |
|----------------|---------------|--------|--|--|--|--|
| 12 HOUR STD    | 467640        | 11.172 |  |  |  |  |
| UPPER LIMIT    | 935280        | 11.672 |  |  |  |  |
| LOWER LIMIT    | 233820        | 10.672 |  |  |  |  |
| EPA SAMPLE NO. |               |        |  |  |  |  |
| MW-1           | 308059        | 11.18  |  |  |  |  |
| MW-2           | 261173        | 11.17  |  |  |  |  |
| MW-3           | 400173        | 11.17  |  |  |  |  |
| MW-4           | 336486        | 11.17  |  |  |  |  |
| TRIP-BLANK     | 279370        | 11.18  |  |  |  |  |
| VV1003WBL01    | 433083        | 11.17  |  |  |  |  |
| VV1003WBS01    | 506399        | 11.17  |  |  |  |  |
| VV1003WBSD01   | 481201        | 11.17  |  |  |  |  |

IS4 = 1,4-Dichlorobenzene-d4

AREA UPPER LIMIT = +100% of internal standard area

AREA LOWER LIMIT = -50% of internal standard area

RT UPPER LIMIT = +0.50 minutes of internal standard RT

RT LOWER LIMIT = -0.50 minutes of internal standard RT

# Column used to flag values outside QC limits with an asterisk.

\* Values outside of QC limits.



# QC SAMPLE DATA

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBL01                        |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBL01                        |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039244.D        | 1         | 10/03/25 11:26 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 1.00  | U         | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 1.00  | U         | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 1.00  | U         | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 5.00  | U         | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 5.00  | U         | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 1.00  | U         | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 1.00  | U         | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 5.00  | U         | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 5.00  | U         | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 1.00  | U         | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 1.00  | U         | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 1.00  | U         | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 5.00  | U         | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 1.00  | U         | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 1.00  | U         | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 1.00  | U         | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 5.00  | U         | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 1.00  | U         | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 1.00  | U         | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 1.00  | U         | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 1.00  | U         | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 1.00  | U         | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 1.00  | U         | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBL01                        |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBL01                        |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039244.D        | 1         | 10/03/25 11:26 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 3.00   | U         | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 1.00   | U         | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 1.00   | U         | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 1.00   | U         | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 1.00   | U         | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 1.00   | U         | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 1.00   | U         | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 5.00   | U         | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 5.00   | U         | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 51.7   |           | 74 - 125 | 103%       | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 52.5   |           | 75 - 124 | 105%       | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 44.6   |           | 86 - 113 | 89%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 45.2   |           | 77 - 121 | 90%        | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 502000 | 4.6       |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 854000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 854000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 433000 | 11.172    |          |            |         |

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

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBS01                        |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBS01                        |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039245.D        | 1         | 10/03/25 13:07 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 17.5  |           | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 18.1  |           | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 19.1  |           | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 18.8  |           | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 88.9  |           | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 91.0  |           | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 16.0  |           | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 19.9  |           | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 88.0  |           | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 89.1  |           | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 19.2  |           | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 19.2  |           | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 19.3  |           | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 18.1  |           | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 18.2  |           | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 19.4  |           | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 19.9  |           | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 19.8  |           | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 19.3  |           | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 100   |           | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 20.9  |           | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 21.1  |           | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 20.9  |           | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 91.8  |           | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 20.7  |           | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 20.6  |           | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 20.3  |           | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 20.2  |           | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 20.3  |           | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 20.2  |           | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBS01                        |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBS01                        |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039245.D        | 1         | 10/03/25 13:07 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 64.6   |           | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 19.7   |           | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 21.5   |           | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 18.2   |           | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 18.1   |           | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 19.3   |           | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 19.7   |           | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 20.5   |           | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 22.0   |           | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 19.0   |           | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 41.7   |           | 74 - 125 | 83%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 49.6   |           | 75 - 124 | 99%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.8   |           | 86 - 113 | 92%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 53.4   |           | 77 - 121 | 107%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 559000 | 4.596     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 866000 | 5.532     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 845000 | 8.773     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 506000 | 11.172    |          |            |         |

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

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBSD01                       |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBSD01                       |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039246.D        | 1         | 10/03/25 13:29 | VV100325      |

| CAS Number     | Parameter                 | Conc. | Qualifier | MDL  | LOQ / CRQL | Units |
|----------------|---------------------------|-------|-----------|------|------------|-------|
| <b>TARGETS</b> |                           |       |           |      |            |       |
| 75-01-4        | Vinyl Chloride            | 17.0  |           | 0.26 | 1.00       | ug/L  |
| 75-00-3        | Chloroethane              | 17.8  |           | 0.47 | 1.00       | ug/L  |
| 75-69-4        | Trichlorofluoromethane    | 18.8  |           | 0.33 | 1.00       | ug/L  |
| 75-35-4        | 1,1-Dichloroethene        | 18.7  |           | 0.23 | 1.00       | ug/L  |
| 107-13-1       | Acrylonitrile             | 94.7  |           | 0.83 | 5.00       | ug/L  |
| 67-64-1        | Acetone                   | 93.4  |           | 1.50 | 5.00       | ug/L  |
| 75-15-0        | Carbon Disulfide          | 15.9  |           | 0.21 | 1.00       | ug/L  |
| 75-09-2        | Methylene Chloride        | 20.2  |           | 0.28 | 1.00       | ug/L  |
| 108-05-4       | Vinyl Acetate             | 92.0  |           | 0.66 | 5.00       | ug/L  |
| 78-93-3        | 2-Butanone                | 93.5  |           | 0.98 | 5.00       | ug/L  |
| 56-23-5        | Carbon Tetrachloride      | 18.8  |           | 0.25 | 1.00       | ug/L  |
| 156-59-2       | cis-1,2-Dichloroethene    | 19.6  |           | 0.19 | 1.00       | ug/L  |
| 74-97-5        | Bromochloromethane        | 19.6  |           | 0.22 | 1.00       | ug/L  |
| 67-66-3        | Chloroform                | 18.2  |           | 0.25 | 1.00       | ug/L  |
| 71-55-6        | 1,1,1-Trichloroethane     | 18.4  |           | 0.20 | 1.00       | ug/L  |
| 71-43-2        | Benzene                   | 19.3  |           | 0.15 | 1.00       | ug/L  |
| 78-87-5        | 1,2-Dichloropropane       | 19.1  |           | 0.20 | 1.00       | ug/L  |
| 74-95-3        | Dibromomethane            | 20.1  |           | 0.25 | 1.00       | ug/L  |
| 75-27-4        | Bromodichloromethane      | 19.4  |           | 0.22 | 1.00       | ug/L  |
| 108-10-1       | 4-Methyl-2-Pentanone      | 110   |           | 0.68 | 5.00       | ug/L  |
| 108-88-3       | Toluene                   | 20.9  |           | 0.14 | 1.00       | ug/L  |
| 10061-02-6     | t-1,3-Dichloropropene     | 21.3  |           | 0.17 | 1.00       | ug/L  |
| 10061-01-5     | cis-1,3-Dichloropropene   | 21.1  |           | 0.16 | 1.00       | ug/L  |
| 591-78-6       | 2-Hexanone                | 95.2  |           | 0.89 | 5.00       | ug/L  |
| 124-48-1       | Dibromochloromethane      | 20.6  |           | 0.18 | 1.00       | ug/L  |
| 106-93-4       | 1,2-Dibromoethane         | 21.4  |           | 0.15 | 1.00       | ug/L  |
| 127-18-4       | Tetrachloroethene         | 20.4  |           | 0.23 | 1.00       | ug/L  |
| 108-90-7       | Chlorobenzene             | 20.2  |           | 0.12 | 1.00       | ug/L  |
| 630-20-6       | 1,1,1,2-Tetrachloroethane | 20.7  |           | 0.19 | 1.00       | ug/L  |
| 100-41-4       | Ethyl Benzene             | 20.2  |           | 0.13 | 1.00       | ug/L  |

## Report of Analysis

|                    |                                    |           |                 |              |
|--------------------|------------------------------------|-----------|-----------------|--------------|
| Client:            | Lockwood, Kessler & Bartlett, Inc. |           | Date Collected: |              |
| Project:           | Ansonia Landfill 2025              |           | Date Received:  |              |
| Client Sample ID:  | VV1003WBSD01                       |           | SDG No.:        | Q3254        |
| Lab Sample ID:     | VV1003WBSD01                       |           | Matrix:         | Water        |
| Analytical Method: | 8260D                              |           | % Solid:        | 0            |
| Sample Wt/Vol:     | 5                                  | Units: mL | Final Vol:      | 5000 uL      |
| Soil Aliquot Vol:  |                                    | uL        | Test:           | VOCMS Group1 |
| GC Column:         | DB-624UI                           | ID : 0.18 | Level :         | LOW          |
| Prep Method :      |                                    |           |                 |              |

|                   |           |                |               |
|-------------------|-----------|----------------|---------------|
| File ID/Qc Batch: | Dilution: | Date Analyzed  | Prep Batch ID |
| VV039246.D        | 1         | 10/03/25 13:29 | VV100325      |

| CAS Number                | Parameter                   | Conc.  | Qualifier | MDL      | LOQ / CRQL | Units   |
|---------------------------|-----------------------------|--------|-----------|----------|------------|---------|
| 1330-20-7                 | Total Xylenes               | 65.1   |           | 0.36     | 3.00       | ug/L    |
| 100-42-5                  | Styrene                     | 19.6   |           | 0.15     | 1.00       | ug/L    |
| 75-25-2                   | Bromoform                   | 22.2   |           | 0.19     | 1.00       | ug/L    |
| 79-34-5                   | 1,1,2,2-Tetrachloroethane   | 18.8   |           | 0.26     | 1.00       | ug/L    |
| 96-18-4                   | 1,2,3-Trichloropropane      | 18.8   |           | 0.35     | 1.00       | ug/L    |
| 106-46-7                  | 1,4-Dichlorobenzene         | 19.5   |           | 0.19     | 1.00       | ug/L    |
| 95-50-1                   | 1,2-Dichlorobenzene         | 20.2   |           | 0.16     | 1.00       | ug/L    |
| 96-12-8                   | 1,2-Dibromo-3-Chloropropane | 20.6   |           | 0.53     | 1.00       | ug/L    |
| 74-88-4                   | Methyl Iodide               | 22.2   |           | 0.83     | 5.00       | ug/L    |
| 110-57-6                  | trans-1,4-Dichloro-2-butene | 19.5   |           | 0.84     | 5.00       | ug/L    |
| <b>SURROGATES</b>         |                             |        |           |          |            |         |
| 17060-07-0                | 1,2-Dichloroethane-d4       | 42.8   |           | 74 - 125 | 86%        | SPK: 50 |
| 1868-53-7                 | Dibromofluoromethane        | 48.6   |           | 75 - 124 | 97%        | SPK: 50 |
| 2037-26-5                 | Toluene-d8                  | 45.3   |           | 86 - 113 | 91%        | SPK: 50 |
| 460-00-4                  | 4-Bromofluorobenzene        | 53.2   |           | 77 - 121 | 106%       | SPK: 50 |
| <b>INTERNAL STANDARDS</b> |                             |        |           |          |            |         |
| 363-72-4                  | Pentafluorobenzene          | 531000 | 4.603     |          |            |         |
| 540-36-3                  | 1,4-Difluorobenzene         | 835000 | 5.535     |          |            |         |
| 3114-55-4                 | Chlorobenzene-d5            | 808000 | 8.776     |          |            |         |
| 3855-82-1                 | 1,4-Dichlorobenzene-d4      | 481000 | 11.172    |          |            |         |

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



# CALIBRATION SUMMARY

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                     |                                      |                      |                                     |
|---------------------|--------------------------------------|----------------------|-------------------------------------|
| Lab Name:           | <u>Alliance</u>                      | Contract:            | <u>LOCK01</u>                       |
| Lab Code:           | <u>ACE</u>                           | SDG No.:             | <u>Q3254</u>                        |
| Instrument ID:      | <u>MSVOA_V</u>                       | Calibration Date(s): | <u>09/22/2025</u> <u>09/22/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>                             | Calibration Time(s): | <u>12:26</u> <u>16:35</u>           |
| GC Column:          | <u>DB-624UI</u> ID: <u>0.18</u> (mm) |                      |                                     |

|                           |        |                     |        |                     |        |                     |       |       |
|---------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:              |        | RRF010 = VV039137.D |        | RRF020 = VV039138.D |        | RRF050 = VV039139.D |       |       |
|                           |        | RRF100 = VV039140.D |        | RRF005 = VV039142.D |        | RRF001 = VV039143.D |       |       |
| COMPOUND                  | RRF010 | RRF020              | RRF050 | RRF100              | RRF005 | RRF001              | RRF   | % RSD |
| Vinyl Chloride            | 1.007  | 0.856               | 0.836  | 0.801               | 0.932  | 1.026               | 0.910 | 10.3  |
| Chloroethane              | 0.662  | 0.568               | 0.533  | 0.500               | 0.606  | 0.915               | 0.631 | 23.8  |
| Trichlorofluoromethane    | 1.394  | 1.253               | 1.193  | 1.151               | 1.365  | 1.546               | 1.317 | 11.2  |
| 1,1-Dichloroethene        | 0.804  | 0.721               | 0.674  | 0.663               | 0.789  | 0.873               | 0.754 | 10.9  |
| Acrylonitrile             | 0.507  | 0.462               | 0.436  | 0.427               | 0.522  | 0.587               | 0.490 | 12.4  |
| Acetone                   | 0.537  | 0.458               | 0.449  | 0.414               | 0.569  | 0.712               | 0.523 | 20.8  |
| Carbon Disulfide          | 2.453  | 2.181               | 2.081  | 1.986               | 2.462  | 2.670               | 2.306 | 11.4  |
| Methylene Chloride        | 0.930  | 0.848               | 0.779  | 0.742               | 1.113  | 1.987               | 1.066 | 44    |
| Vinyl Acetate             | 2.393  | 2.241               | 2.204  | 2.128               | 2.313  | 2.243               | 2.254 | 4     |
| 2-Butanone                | 0.604  | 0.584               | 0.586  | 0.580               | 0.570  | 0.554               | 0.579 | 2.9   |
| Carbon Tetrachloride      | 0.742  | 0.710               | 0.703  | 0.675               | 0.736  | 0.890               | 0.743 | 10.2  |
| cis-1,2-Dichloroethene    | 0.908  | 0.842               | 0.868  | 0.860               | 0.832  | 1.001               | 0.885 | 7.1   |
| Bromochloromethane        | 0.836  | 0.753               | 0.679  | 0.637               | 0.525  | 0.743               | 0.696 | 15.5  |
| Chloroform                | 1.752  | 1.570               | 1.490  | 1.433               | 1.746  | 1.880               | 1.645 | 10.6  |
| 1,1,1-Trichloroethane     | 1.562  | 1.321               | 1.298  | 1.263               | 1.513  | 1.590               | 1.424 | 10.3  |
| Benzene                   | 2.077  | 1.978               | 1.990  | 1.931               | 1.963  | 1.936               | 1.979 | 2.7   |
| 1,2-Dichloropropane       | 0.509  | 0.519               | 0.502  | 0.486               | 0.471  | 0.465               | 0.492 | 4.4   |
| Dibromomethane            | 0.392  | 0.367               | 0.368  | 0.354               | 0.381  | 0.396               | 0.376 | 4.4   |
| Bromodichloromethane      | 0.832  | 0.764               | 0.753  | 0.725               | 0.767  | 0.829               | 0.778 | 5.5   |
| 4-Methyl-2-Pentanone      | 0.715  | 0.736               | 0.732  | 0.714               | 0.613  | 0.459               | 0.661 | 16.5  |
| Toluene                   | 1.211  | 1.199               | 1.229  | 1.214               | 1.130  | 0.901               | 1.147 | 10.9  |
| t-1,3-Dichloropropene     | 0.701  | 0.696               | 0.728  | 0.753               | 0.645  | 0.548               | 0.678 | 10.8  |
| cis-1,3-Dichloropropene   | 0.803  | 0.777               | 0.811  | 0.803               | 0.704  | 0.595               | 0.749 | 11.4  |
| 2-Hexanone                | 0.505  | 0.534               | 0.546  | 0.533               | 0.429  | 0.231               | 0.463 | 26.2  |
| Dibromochloromethane      | 0.605  | 0.570               | 0.567  | 0.558               | 0.598  | 0.620               | 0.586 | 4.2   |
| 1,2-Dibromoethane         | 0.556  | 0.538               | 0.529  | 0.514               | 0.535  | 0.484               | 0.526 | 4.7   |
| Tetrachloroethene         | 0.435  | 0.412               | 0.412  | 0.407               | 0.431  | 0.415               | 0.419 | 2.7   |
| Chlorobenzene             | 1.403  | 1.368               | 1.383  | 1.358               | 1.417  | 1.428               | 1.393 | 2     |
| 1,1,1,2-Tetrachloroethane | 0.520  | 0.488               | 0.488  | 0.477               | 0.504  | 0.543               | 0.503 | 4.8   |
| Ethyl Benzene             | 2.028  | 2.143               | 2.382  | 2.418               | 1.953  | 1.725               | 2.108 | 12.5  |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                                                 |                                                          |
|-------------------------------------------------|----------------------------------------------------------|
| Lab Name: <u>Alliance</u>                       | Contract: <u>LOCK01</u>                                  |
| Lab Code: <u>ACE</u>                            | SDG No.: <u>Q3254</u>                                    |
| Instrument ID: <u>MSVOA_V</u>                   | Calibration Date(s): <u>09/22/2025</u> <u>09/22/2025</u> |
| Heated Purge: (Y/N) <u>N</u>                    | Calibration Time(s): <u>12:26</u> <u>16:35</u>           |
| GC Column: <u>DB-624UI</u> ID: <u>0.18</u> (mm) |                                                          |

|                             |        |                     |        |                     |        |                     |       |       |
|-----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                |        | RRF010 = VV039137.D |        | RRF020 = VV039138.D |        | RRF050 = VV039139.D |       |       |
|                             |        | RRF100 = VV039140.D |        | RRF005 = VV039142.D |        | RRF001 = VV039143.D |       |       |
| COMPOUND                    | RRF010 | RRF020              | RRF050 | RRF100              | RRF005 | RRF001              | RRF   | % RSD |
| m/p-Xylenes                 | 0.820  | 0.882               | 0.949  | 0.938               | 0.713  | 0.580               | 0.814 | 17.7  |
| o-Xylene                    | 0.703  | 0.742               | 0.846  | 0.880               | 0.635  | 0.518               | 0.720 | 18.7  |
| Styrene                     | 1.239  | 1.381               | 1.553  | 1.537               | 1.038  | 0.828               | 1.263 | 22.8  |
| Bromoform                   | 0.396  | 0.384               | 0.396  | 0.400               | 0.385  | 0.397               | 0.393 | 1.7   |
| 1,1,2,2-Tetrachloroethane   | 1.600  | 1.461               | 1.425  | 1.389               | 1.684  | 1.975               | 1.589 | 13.8  |
| 1,2,3-Trichloropropane      | 1.136  | 1.048               | 1.049  | 1.023               | 1.201  | 1.375               | 1.138 | 11.7  |
| 1,4-Dichlorobenzene         | 2.070  | 1.951               | 1.934  | 1.919               | 2.139  | 2.522               | 2.089 | 11    |
| 1,2-Dichlorobenzene         | 1.751  | 1.652               | 1.738  | 1.802               | 1.797  | 1.935               | 1.779 | 5.3   |
| 1,2-Dibromo-3-Chloropropane | 0.299  | 0.283               | 0.285  | 0.302               | 0.333  | 0.476               | 0.330 | 22.4  |
| 1,2-Dichloroethane-d4       | 0.773  | 0.663               | 0.646  | 0.618               | 0.919  |                     | 0.724 | 17.1  |
| Dibromofluoromethane        | 0.435  | 0.382               | 0.377  | 0.360               | 0.454  |                     | 0.402 | 10.1  |
| Toluene-d8                  | 1.166  | 1.092               | 1.136  | 1.097               | 1.411  |                     | 1.181 | 11.2  |
| 4-Bromofluorobenzene        | 0.481  | 0.464               | 0.507  | 0.501               | 0.477  |                     | 0.486 | 3.6   |
| Methyl Iodide               | 0.772  | 0.742               | 0.786  | 0.793               | 0.638  |                     | 0.746 | 8.5   |
| trans-1,4-Dichloro-2-butene | 0.478  | 0.451               | 0.495  | 0.521               | 0.427  |                     | 0.474 | 7.8   |

\* Compounds with required minimum RRF and maximum %RSD values.  
All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                              |
|---------------------|-------------------|------------------------|------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>LOCK01</u>                |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q3254</u>                 |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>10/03/2025 08:53</u>      |
| Lab File ID:        | <u>VV039242.D</u> | Init. Calib. Date(s):  | <u>09/22/2025 09/22/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>12:26 16:35</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                  | RRF   | RRF050 | MIN RRF | %D     | MAX%D |
|---------------------------|-------|--------|---------|--------|-------|
| Vinyl Chloride            | 0.910 | 0.797  |         | -12.42 | 20    |
| Chloroethane              | 0.631 | 0.533  |         | -15.53 | 20    |
| Trichlorofluoromethane    | 1.317 | 1.296  |         | -1.6   | 20    |
| 1,1-Dichloroethene        | 0.754 | 0.737  |         | -2.26  | 20    |
| Acrylonitrile             | 0.490 | 0.450  |         | -8.16  | 20    |
| Acetone                   | 0.523 | 0.465  |         | -11.09 | 20    |
| Carbon Disulfide          | 2.306 | 1.914  |         | -17    | 20    |
| Methylene Chloride        | 1.066 | 0.842  |         | -21.01 | 20    |
| Vinyl Acetate             | 2.254 | 2.045  |         | -9.27  | 20    |
| 2-Butanone                | 0.579 | 0.545  |         | -5.87  | 20    |
| Carbon Tetrachloride      | 0.743 | 0.816  |         | 9.82   | 20    |
| cis-1,2-Dichloroethene    | 0.885 | 0.850  |         | -3.95  | 20    |
| Bromochloromethane        | 0.696 | 0.695  |         | -0.14  | 20    |
| Chloroform                | 1.645 | 1.541  |         | -6.32  | 20    |
| 1,1,1-Trichloroethane     | 1.424 | 1.326  |         | -6.88  | 20    |
| Benzene                   | 1.979 | 2.157  |         | 8.99   | 20    |
| 1,2-Dichloropropane       | 0.492 | 0.557  |         | 13.21  | 20    |
| Dibromomethane            | 0.376 | 0.419  |         | 11.44  | 20    |
| Bromodichloromethane      | 0.778 | 0.870  |         | 11.82  | 20    |
| 4-Methyl-2-Pentanone      | 0.661 | 0.770  |         | 16.49  | 20    |
| Toluene                   | 1.147 | 1.390  |         | 21.19  | 20    |
| t-1,3-Dichloropropene     | 0.678 | 0.811  |         | 19.62  | 20    |
| cis-1,3-Dichloropropene   | 0.749 | 0.891  |         | 18.96  | 20    |
| 2-Hexanone                | 0.463 | 0.579  |         | 25.05  | 20    |
| Dibromochloromethane      | 0.586 | 0.698  |         | 19.11  | 20    |
| 1,2-Dibromoethane         | 0.526 | 0.621  |         | 18.06  | 20    |
| Tetrachloroethene         | 0.419 | 0.472  |         | 12.65  | 20    |
| Chlorobenzene             | 1.393 | 1.547  | 0.3     | 10.98  | 20    |
| 1,1,1,2-Tetrachloroethane | 0.503 | 0.582  |         | 15.71  | 20    |
| Ethyl Benzene             | 2.108 | 2.531  |         | 20.07  | 20    |
| m/p-Xylenes               | 0.814 | 1.051  |         | 29.11  | 20    |
| o-Xylene                  | 0.720 | 0.926  |         | 28.61  | 20    |
| Styrene                   | 1.263 | 1.770  |         | 40.14  | 20    |
| Bromoform                 | 0.393 | 0.490  | 0.1     | 24.68  | 20    |
| 1,1,2,2-Tetrachloroethane | 1.589 | 1.472  | 0.3     | -7.36  | 20    |
| 1,2,3-Trichloropropane    | 1.138 | 1.053  |         | -7.47  | 20    |
| 1,4-Dichlorobenzene       | 2.089 | 2.097  |         | 0.38   | 20    |
| 1,2-Dichlorobenzene       | 1.779 | 1.851  |         | 4.05   | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                              |
|---------------------|-------------------|------------------------|------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>LOCK01</u>                |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q3254</u>                 |
| Instrument ID:      | <u>MSVOA_V</u>    | Calibration Date/Time: | <u>10/03/2025 08:53</u>      |
| Lab File ID:        | <u>VV039242.D</u> | Init. Calib. Date(s):  | <u>09/22/2025 09/22/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>          | Init. Calib. Time(s):  | <u>12:26 16:35</u>           |
| GC Column:          | <u>DB-624UI</u>   | ID:                    | <u>0.18</u> (mm)             |

| COMPOUND                    | RRF   | RRF050 | MIN<br>RRF | %D     | MAX%D |
|-----------------------------|-------|--------|------------|--------|-------|
| 1,2-Dibromo-3-Chloropropane | 0.330 | 0.287  |            | -13.03 | 20    |
| 1,2-Dichloroethane-d4       | 0.724 | 0.604  |            | -16.58 | 20    |
| Dibromofluoromethane        | 0.402 | 0.442  |            | 9.95   | 20    |
| Toluene-d8                  | 1.181 | 1.194  |            | 1.1    | 20    |
| 4-Bromofluorobenzene        | 0.486 | 0.586  |            | 20.58  | 20    |
| Methyl Iodide               | 0.746 | 0.854  |            | 14.48  | 20    |
| trans-1,4-Dichloro-2-butene | 0.474 | 0.497  |            | 4.85   | 20    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.