

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: ENTA05  
 Lab Code: CHEM Case No.: Q1282 SAS No.: Q1282 SDG No.: Q1282  
 Instrument ID: MSVOA\_N Calibration Date(s): 01/14/2025 01/14/2025  
 Heated Purge: (Y/N) N Calibration Time(s): 14:56 17:19  
 GC Column: RXI-624 ID: 0.25 (mm)

| LAB FILE ID:            |        | RRF100 = VN085438.D |        | RRF050 = VN085439.D |        | RRF020 = VN085440.D |       |       |  |
|-------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                         |        | RRF010 = VN085441.D |        | RRF005 = VN085442.D |        | RRF001 = VN085443.D |       |       |  |
| COMPOUND                | RRF100 | RRF050              | RRF020 | RRF010              | RRF005 | RRF001              | RRF   | % RSD |  |
| Methyl tert-butyl Ether | 1.834  | 1.873               | 1.853  | 1.664               | 1.685  | 1.545               | 1.742 | 7.5   |  |
| Carbon Tetrachloride    | 0.574  | 0.530               | 0.579  | 0.529               | 0.565  | 0.567               | 0.557 | 4     |  |
| Chloroform              | 1.197  | 1.175               | 1.241  | 1.169               | 1.253  | 1.273               | 1.218 | 3.6   |  |
| 1,1,1-Trichloroethane   | 1.053  | 1.016               | 1.091  | 1.000               | 1.148  | 1.102               | 1.068 | 5.2   |  |
| Benzene                 | 1.551  | 1.449               | 1.527  | 1.376               | 1.474  | 1.400               | 1.463 | 4.7   |  |
| Toluene                 | 0.964  | 0.870               | 0.919  | 0.808               | 0.835  | 0.690               | 0.848 | 11.3  |  |
| Tetrachloroethene       | 0.351  | 0.322               | 0.365  | 0.338               | 0.346  | 0.323               | 0.341 | 4.9   |  |
| Ethyl Benzene           | 2.072  | 1.867               | 1.940  | 1.685               | 1.709  | 1.430               | 1.784 | 12.7  |  |
| m/p-Xylenes             | 0.775  | 0.707               | 0.750  | 0.615               | 0.616  | 0.492               | 0.659 | 16    |  |
| o-Xylene                | 0.738  | 0.681               | 0.713  | 0.584               | 0.582  | 0.482               | 0.630 | 15.5  |  |
| 1,2-Dichloroethane-d4   | 0.774  | 0.831               | 0.754  | 0.762               | 0.914  |                     | 0.807 | 8.3   |  |
| Dibromofluoromethane    | 0.359  | 0.358               | 0.335  | 0.310               | 0.373  |                     | 0.347 | 7.1   |  |
| Toluene-d8              | 1.339  | 1.267               | 1.207  | 1.076               | 1.274  |                     | 1.232 | 8.1   |  |
| 4-Bromofluorobenzene    | 0.475  | 0.449               | 0.410  | 0.357               | 0.417  |                     | 0.422 | 10.6  |  |

\* 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.