

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: RUTW01  
 Lab Code: CHEM Case No.: Q1241 SAS No.: Q1241 SDG No.: Q1241  
 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 |  |
| Vinyl Chloride        | 0.697  | 0.686               | 0.727  | 0.711               | 0.781  | 0.819               | 0.737 | 7.1   |  |
| 1,1-Dichloroethene    | 0.548  | 0.533               | 0.556  | 0.526               | 0.559  | 0.497               | 0.537 | 4.3   |  |
| 2-Butanone            | 0.378  | 0.390               | 0.398  | 0.363               | 0.387  | 0.386               | 0.384 | 3.1   |  |
| 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   |  |
| Benzene               | 1.551  | 1.449               | 1.527  | 1.376               | 1.474  | 1.400               | 1.463 | 4.7   |  |
| 1,2-Dichloroethane    | 0.569  | 0.547               | 0.575  | 0.522               | 0.574  | 0.517               | 0.551 | 4.8   |  |
| Trichloroethene       | 0.362  | 0.324               | 0.352  | 0.310               | 0.343  | 0.352               | 0.341 | 5.8   |  |
| Tetrachloroethene     | 0.351  | 0.322               | 0.365  | 0.338               | 0.346  | 0.323               | 0.341 | 4.9   |  |
| Chlorobenzene         | 1.133  | 1.076               | 1.154  | 1.047               | 1.110  | 1.051               | 1.095 | 4     |  |
| 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.