

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: JAC005  
 Lab Code: CHEM Case No.: Q2234 SAS No.: Q2234 SDG No.: Q2234  
 Instrument ID: MSVOA\_X Calibration Date(s): 06/06/2025 06/06/2025  
 Heated Purge: (Y/N) N Calibration Time(s): 09:42 12:57  
 GC Column: DB-624UI ID: 0.18 (mm)

| LAB FILE ID:           |        | RRF005 = VX046518.D |        | RRF020 = VX046519.D |        | RRF050 = VX046520.D |       |       |  |
|------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                        |        | RRF100 = VX046521.D |        | RRF150 = VX046522.D |        | RRF001 = VX046524.D |       |       |  |
| COMPOUND               | RRF005 | RRF020              | RRF050 | RRF100              | RRF150 | RRF001              | RRF   | % RSD |  |
| Vinyl Chloride         | 0.697  | 0.593               | 0.596  | 0.591               | 0.622  | 0.679               | 0.630 | 7.5   |  |
| 1,1-Dichloroethene     | 0.663  | 0.550               | 0.567  | 0.561               | 0.585  | 0.635               | 0.594 | 7.6   |  |
| 1,1-Dichloroethane     | 1.349  | 1.266               | 1.259  | 1.234               | 1.297  | 1.281               | 1.281 | 3.1   |  |
| cis-1,2-Dichloroethene | 0.786  | 0.752               | 0.741  | 0.728               | 0.767  | 0.866               | 0.773 | 6.4   |  |
| 1,1,1-Trichloroethane  | 1.170  | 1.131               | 1.141  | 1.128               | 1.188  | 1.131               | 1.148 | 2.2   |  |
| Benzene                | 1.597  | 1.503               | 1.427  | 1.380               | 1.442  | 1.522               | 1.479 | 5.3   |  |
| 1,2-Dichloroethane     | 0.646  | 0.641               | 0.610  | 0.586               | 0.602  | 0.606               | 0.615 | 3.8   |  |
| Trichloroethene        | 0.385  | 0.356               | 0.351  | 0.332               | 0.354  | 0.476               | 0.376 | 13.8  |  |
| 1,1,2-Trichloroethane  | 0.375  | 0.389               | 0.366  | 0.356               | 0.372  | 0.331               | 0.365 | 5.4   |  |
| Tetrachloroethene      | 0.347  | 0.324               | 0.310  | 0.301               | 0.314  | 0.410               | 0.334 | 12.1  |  |
| 1,2-Dichloroethane-d4  | 0.997  | 0.848               | 0.873  | 0.828               | 0.900  |                     | 0.890 | 7.4   |  |
| Dibromofluoromethane   | 0.392  | 0.353               | 0.368  | 0.347               | 0.379  |                     | 0.368 | 5     |  |
| Toluene-d8             | 1.362  | 1.159               | 1.188  | 1.132               | 1.220  |                     | 1.212 | 7.4   |  |
| 4-Bromofluorobenzene   | 0.564  | 0.482               | 0.493  | 0.468               | 0.501  |                     | 0.502 | 7.4   |  |

\* 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.