

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: CLOU03  
Lab Code: CHEM Case No.: P4960 SAS No.: P4960 SDG No.: P4960  
Instrument ID: MSVOA\_X Calibration Date(s): 11/21/2024 11/21/2024  
Heated Purge: (Y/N) N Calibration Time(s): 09:47 12:03  
GC Column: DB-624UI ID: 0.18 (mm)

| LAB FILE ID:            |        | RRF001 = VX043925.D |        | RRF005 = VX043926.D |        | RRF020 = VX043927.D |       |       |  |
|-------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                         |        | RRF050 = VX043928.D |        | RRF100 = VX043929.D |        | RRF150 = VX043930.D |       |       |  |
| COMPOUND                | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |  |
| Methyl tert-butyl Ether | 1.732  | 1.991               | 2.212  | 2.264               | 2.160  | 2.192               | 2.092 | 9.5   |  |
| Benzene                 | 1.211  | 1.369               | 1.457  | 1.494               | 1.388  | 1.402               | 1.387 | 7     |  |
| Toluene                 | 0.636  | 0.839               | 0.883  | 0.921               | 0.850  | 0.852               | 0.830 | 12    |  |
| Ethyl Benzene           | 1.579  | 1.831               | 1.910  | 2.024               | 1.884  | 1.938               | 1.861 | 8.2   |  |
| m/p-Xylenes             | 0.610  | 0.663               | 0.704  | 0.762               | 0.700  | 0.725               | 0.694 | 7.6   |  |
| o-Xylene                | 0.543  | 0.665               | 0.691  | 0.757               | 0.693  | 0.716               | 0.678 | 10.7  |  |
| Isopropylbenzene        | 3.435  | 3.927               | 3.996  | 4.097               | 3.679  | 3.927               | 3.843 | 6.3   |  |
| n-propylbenzene         | 3.667  | 4.115               | 4.469  | 4.733               | 4.286  | 4.486               | 4.293 | 8.6   |  |
| 1,3,5-Trimethylbenzene  | 2.704  | 3.115               | 3.352  | 3.444               | 3.118  | 3.238               | 3.162 | 8.2   |  |
| tert-Butylbenzene       | 2.581  | 3.097               | 3.266  | 3.440               | 3.152  | 3.269               | 3.134 | 9.4   |  |
| 1,2,4-Trimethylbenzene  | 2.557  | 3.173               | 3.366  | 3.440               | 3.124  | 3.228               | 3.148 | 9.9   |  |
| sec-Butylbenzene        | 3.116  | 3.921               | 3.968  | 4.249               | 3.812  | 3.974               | 3.840 | 10    |  |
| p-Isopropyltoluene      | 2.255  | 3.043               | 3.316  | 3.521               | 3.219  | 3.362               | 3.119 | 14.5  |  |
| n-Butylbenzene          | 1.897  | 2.580               | 2.805  | 3.184               | 2.909  | 3.126               | 2.750 | 17.2  |  |
| Naphthalene             | 2.605  | 3.361               | 3.698  | 4.005               | 3.733  | 4.058               | 3.576 | 15    |  |
| 1,2-Dichloroethane-d4   |        | 0.926               | 0.876  | 0.751               | 0.855  | 0.880               | 0.858 | 7.6   |  |
| Dibromofluoromethane    |        | 0.371               | 0.353  | 0.327               | 0.359  | 0.376               | 0.357 | 5.4   |  |
| Toluene-d8              |        | 1.327               | 1.237  | 1.104               | 1.232  | 1.257               | 1.232 | 6.6   |  |
| 4-Bromofluorobenzene    |        | 0.401               | 0.414  | 0.387               | 0.438  | 0.455               | 0.419 | 6.6   |  |

\* 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.