

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: TETR06  
Lab Code: CHEM Case No.: P5316 SAS No.: P5316 SDG No.: P5316  
Instrument ID: MSVOA\_Y Calibration Date(s): 12/26/2024 12/26/2024  
Heated Purge: (Y/N) Y Calibration Time(s): 09:31 11:51  
GC Column: RXI-624 ID: 0.25 (mm)

|                          |                     |        |                     |        |                     |        |       |       |
|--------------------------|---------------------|--------|---------------------|--------|---------------------|--------|-------|-------|
| LAB FILE ID:             | RRF005 = VY020706.D |        | RRF010 = VY020707.D |        | RRF020 = VY020708.D |        |       |       |
|                          | RRF050 = VY020709.D |        | RRF150 = VY020710.D |        | RRF100 = VY020711.D |        |       |       |
| COMPOUND                 | RRF005              | RRF010 | RRF020              | RRF050 | RRF150              | RRF100 | RRF   | % RSD |
| Vinyl Chloride           | 0.168               | 0.195  | 0.176               | 0.192  | 0.178               | 0.181  | 0.182 | 5.6   |
| 1,1-Dichloroethene       | 0.326               | 0.359  | 0.330               | 0.352  | 0.340               | 0.344  | 0.342 | 3.7   |
| Acetone                  | 0.050               | 0.068  | 0.045               | 0.054  | 0.048               | 0.046  | 0.052 | 16.8  |
| Methyl tert-butyl Ether  | 0.880               | 1.004  | 0.922               | 1.015  | 0.942               | 0.951  | 0.952 | 5.3   |
| Methylene Chloride       | 0.324               | 0.365  | 0.328               | 0.343  | 0.329               | 0.338  | 0.338 | 4.5   |
| trans-1,2-Dichloroethene | 0.339               | 0.378  | 0.356               | 0.391  | 0.368               | 0.378  | 0.368 | 5     |
| 1,1-Dichloroethane       | 0.571               | 0.661  | 0.596               | 0.620  | 0.591               | 0.620  | 0.609 | 5.1   |
| 2-Butanone               | 0.069               | 0.069  | 0.066               | 0.080  | 0.071               | 0.067  | 0.070 | 7     |
| Carbon Tetrachloride     | 0.556               | 0.644  | 0.581               | 0.614  | 0.593               | 0.606  | 0.599 | 5     |
| cis-1,2-Dichloroethene   | 0.419               | 0.460  | 0.426               | 0.462  | 0.432               | 0.447  | 0.441 | 4.1   |
| Chloroform               | 0.839               | 0.821  | 0.753               | 0.788  | 0.742               | 0.762  | 0.784 | 5     |
| 1,1,1-Trichloroethane    | 0.814               | 0.897  | 0.826               | 0.859  | 0.816               | 0.831  | 0.840 | 3.8   |
| Benzene                  | 0.969               | 1.104  | 1.040               | 1.089  | 1.041               | 1.083  | 1.054 | 4.7   |
| 1,2-Dichloroethane       | 0.297               | 0.352  | 0.325               | 0.364  | 0.342               | 0.348  | 0.338 | 7     |
| Trichloroethene          | 0.288               | 0.336  | 0.315               | 0.324  | 0.309               | 0.317  | 0.315 | 5.1   |
| Toluene                  | 0.654               | 0.751  | 0.722               | 0.745  | 0.723               | 0.741  | 0.723 | 4.9   |
| Tetrachloroethene        | 0.308               | 0.360  | 0.348               | 0.351  | 0.337               | 0.346  | 0.342 | 5.3   |
| Chlorobenzene            | 0.925               | 1.044  | 0.976               | 1.002  | 0.943               | 0.985  | 0.979 | 4.3   |
| Ethyl Benzene            | 1.540               | 1.810  | 1.715               | 1.778  | 1.682               | 1.743  | 1.711 | 5.6   |
| m/p-Xylenes              | 0.590               | 0.691  | 0.649               | 0.675  | 0.643               | 0.667  | 0.653 | 5.4   |
| o-Xylene                 | 0.552               | 0.666  | 0.639               | 0.650  | 0.603               | 0.632  | 0.624 | 6.6   |
| n-propylbenzene          | 3.521               | 4.231  | 3.972               | 3.896  | 3.774               | 3.902  | 3.883 | 6     |
| tert-Butylbenzene        | 2.487               | 2.888  | 2.751               | 2.727  | 2.632               | 2.687  | 2.695 | 4.9   |
| 1,2,4-Trimethylbenzene   | 2.513               | 2.946  | 2.860               | 2.888  | 2.777               | 2.867  | 2.809 | 5.5   |
| sec-Butylbenzene         | 3.316               | 3.880  | 3.730               | 3.676  | 3.540               | 3.653  | 3.633 | 5.3   |
| 1,3-Dichlorobenzene      | 1.428               | 1.670  | 1.567               | 1.578  | 1.529               | 1.579  | 1.559 | 5.1   |
| 1,4-Dichlorobenzene      | 1.492               | 1.625  | 1.562               | 1.572  | 1.502               | 1.549  | 1.550 | 3.1   |
| n-Butylbenzene           | 2.356               | 2.894  | 2.764               | 2.757  | 2.683               | 2.760  | 2.702 | 6.8   |
| 1,2-Dichlorobenzene      | 1.197               | 1.374  | 1.362               | 1.410  | 1.328               | 1.347  | 1.336 | 5.5   |
| 1,2-Dichloroethane-d4    | 0.335               | 0.402  | 0.313               | 0.516  | 0.433               | 0.457  | 0.410 | 18.7  |

\* 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.

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| COMPOUND             | RRF005 | RRF010              | RRF020 | RRF050              | RRF150 | RRF100              | $\overline{\text{RRF}}$ | % RSD |
| Dibromofluoromethane | 0.235  | 0.274               | 0.232  | 0.319               | 0.284  | 0.307               | 0.275                   | 13.1  |
| Toluene-d8           | 0.733  | 0.883               | 0.733  | 1.227               | 1.093  | 1.158               | 0.971                   | 22.4  |
| 4-Bromofluorobenzene | 0.303  | 0.378               | 0.319  | 0.426               | 0.373  | 0.397               | 0.366                   | 12.8  |
| 1,4-Dioxane          | 1.014  | 1.439               | 1.316  | 1.680               | 1.562  | 1.428               | 1.407                   | 16.3  |

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