

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                     |                  |                      |                   |
|---------------------|------------------|----------------------|-------------------|
| Lab Name:           | <u>CHEMTECH</u>  | Contract:            | <u>WOOD06</u>     |
| Lab Code:           | <u>CHEM</u>      | Case No.:            | <u>Q1668</u>      |
| Instrument ID:      | <u>MSVOA_L</u>   | SAS No.:             | <u>Q1668</u>      |
| Heated Purge: (Y/N) | <u>N</u>         | SDG No.:             | <u>Q1668</u>      |
| GC Column:          | <u>RTX-1</u>     | Calibration Date(s): | <u>04/17/2025</u> |
| ID:                 | <u>0.32 (mm)</u> | Calibration Time(s): | <u>09:44</u>      |
|                     |                  |                      | <u>12:54</u>      |

|                                |        |                     |        |                     |        |                     |       |       |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                   |        | RRF010 = VL042349.D |        | RRF002 = VL042350.D |        | RRF001 = VL042351.D |       |       |
|                                |        | RRF0.5 = VL042352.D |        | RRF0.1 = VL042353.D |        | RRF.03 = VL042354.D |       |       |
| COMPOUND                       | RRF010 | RRF002              | RRF001 | RRF0.5              | RRF0.1 | RRF.03              | RRF   | % RSD |
| Dichlorodifluoromethane        | 1.658  | 1.827               | 1.837  | 1.831               |        |                     | 1.767 | 5     |
| Chloromethane                  | 0.652  | 0.701               | 0.725  | 0.752               |        |                     | 0.699 | 5.9   |
| Vinyl Chloride                 | 0.638  | 0.689               | 0.709  | 0.731               | 0.911  | 1.002               | 0.761 | 18.4  |
| Bromomethane                   | 0.304  | 0.300               | 0.330  | 0.397               |        |                     | 0.327 | 12.6  |
| Chloroethane                   | 0.253  | 0.280               | 0.287  | 0.303               |        |                     | 0.275 | 7.8   |
| Tetrahydrofuran                | 0.409  | 0.468               | 0.480  | 0.518               |        |                     | 0.461 | 9.2   |
| Trichlorofluoromethane         | 1.567  | 1.746               | 1.806  | 1.684               |        |                     | 1.680 | 6     |
| 1,1,2-Trichlorotrifluoroethane | 1.281  | 1.437               | 1.439  | 1.389               |        |                     | 1.374 | 5.1   |
| Dichlorotetrafluoroethane      | 1.432  | 1.581               | 1.550  | 1.556               |        |                     | 1.517 | 4.3   |
| tert-Butyl alcohol             | 1.804  | 2.103               | 2.132  | 2.263               |        |                     | 2.029 | 9.8   |
| Heptane                        | 1.884  | 2.209               | 2.296  | 2.444               |        |                     | 2.153 | 11.1  |
| 1,1-Dichloroethene             | 0.583  | 0.637               | 0.635  | 0.735               |        |                     | 0.642 | 8.8   |
| Acetone                        | 1.263  | 1.743               | 1.895  | 1.968               |        |                     | 1.637 | 20    |
| Carbon Disulfide               | 1.668  | 1.785               | 1.670  | 1.903               |        |                     | 1.749 | 5.6   |
| Methyl tert-Butyl Ether        | 0.940  | 1.132               | 1.141  | 1.134               |        |                     | 1.070 | 8.6   |
| Methylene Chloride             | 0.498  | 0.582               | 0.728  | 0.875               |        |                     | 0.641 | 24.7  |
| trans-1,2-Dichloroethene       | 0.664  | 0.736               | 0.759  | 0.814               |        |                     | 0.731 | 8.3   |
| 1,1-Dichloroethane             | 1.284  | 1.383               | 1.469  | 1.452               |        |                     | 1.388 | 5.4   |
| Cyclohexane                    | 1.252  | 1.347               | 1.333  | 1.422               |        |                     | 1.320 | 5.5   |
| 2-Butanone                     | 0.687  | 0.818               | 0.826  | 0.886               |        |                     | 0.791 | 9.9   |
| Carbon Tetrachloride           | 0.770  | 0.847               | 0.830  | 0.832               | 0.914  | 1.319               | 0.905 | 20.7  |
| cis-1,2-Dichloroethene         | 1.459  | 1.593               | 1.603  | 1.588               |        |                     | 1.553 | 3.9   |
| Chloroform                     | 2.145  | 2.392               | 2.333  | 2.458               |        |                     | 2.309 | 5.5   |
| 1,1,1-Trichloroethane          | 2.274  | 2.493               | 2.549  | 2.557               | 2.840  | 3.110               | 2.599 | 11    |
| 2,2,4-Trimethylpentane         | 1.708  | 1.962               | 1.916  | 1.935               |        |                     | 1.863 | 5.8   |
| Benzene                        | 1.002  | 1.121               | 1.075  | 1.146               |        |                     | 1.079 | 5.3   |
| 1,2-Dichloroethane             | 0.592  | 0.666               | 0.647  | 0.691               |        |                     | 0.647 | 5.6   |
| Trichloroethene                | 0.395  | 0.447               | 0.448  | 0.443               | 0.505  | 0.643               | 0.471 | 17.6  |
| 1,2-Dichloropropane            | 0.370  | 0.405               | 0.413  | 0.441               |        |                     | 0.402 | 7     |
| Bromodichloromethane           | 0.852  | 0.912               | 0.909  | 0.943               |        |                     | 0.901 | 3.7   |

\* 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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| Lab Code:           | <u>CHEM</u>      | Case No.:            | <u>Q1668</u>      |
| Instrument ID:      | <u>MSVOA_L</u>   | SAS No.:             | <u>Q1668</u>      |
| Heated Purge: (Y/N) | <u>N</u>         | SDG No.:             | <u>Q1668</u>      |
| GC Column:          | <u>RTX-1</u>     | Calibration Date(s): | <u>04/17/2025</u> |
| ID:                 | <u>0.32 (mm)</u> | Calibration Time(s): | <u>09:44</u>      |
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|                           |         | RRF0.5 = VL042352.D |         | RRF0.1 = VL042353.D |        | RRF.03 = VL042354.D |         |       |
| COMPOUND                  | RRF010  | RRF002              | RRF001  | RRF0.5              | RRF0.1 | RRF.03              | RRF     | % RSD |
| 4-Methyl-2-Pentanone      | 1.082   | 1.435               | 1.382   | 1.541               |        |                     | 1.314   | 15.1  |
| Toluene                   | 1.193   | 1.370               | 1.295   | 1.371               |        |                     | 1.294   | 6     |
| t-1,3-Dichloropropene     | 0.459   | 0.509               | 0.510   | 0.514               |        |                     | 0.497   | 4.6   |
| cis-1,3-Dichloropropene   | 0.570   | 0.637               | 0.617   | 0.613               |        |                     | 0.608   | 4.1   |
| 1,1,2-Trichloroethane     | 0.376   | 0.420               | 0.429   | 0.423               |        |                     | 0.410   | 5.3   |
| Dibromochloromethane      | 0.679   | 0.746               | 0.721   | 0.743               |        |                     | 0.723   | 3.7   |
| 1,2-Dibromoethane         | 0.539   | 0.606               | 0.605   | 0.582               | 0.621  |                     | 0.586   | 5.3   |
| Tetrachloroethene         | 0.357   | 0.391               | 0.393   | 0.420               | 0.452  | 0.543               | 0.417   | 15.5  |
| Chlorobenzene             | 0.956   | 1.068               | 1.067   | 1.129               |        |                     | 1.040   | 6.8   |
| Ethyl Benzene             | 1.832   | 2.068               | 2.088   | 2.029               |        |                     | 1.981   | 5.7   |
| m/p-Xylene                | 1.456   | 1.658               | 1.682   | 1.685               |        |                     | 1.596   | 6.9   |
| o-Xylene                  | 1.442   | 1.668               | 1.697   | 1.740               |        |                     | 1.607   | 8.3   |
| Styrene                   | 0.726   | 0.808               | 0.776   | 0.792               |        |                     | 0.767   | 4.7   |
| Bromoform                 | 0.616   | 0.675               | 0.623   | 0.622               |        |                     | 0.633   | 3.8   |
| 1,1,2,2-Tetrachloroethane | 0.873   | 0.968               | 0.985   | 0.990               | 1.176  | 1.407               | 1.042   | 18.1  |
| 2-Chlorotoluene           | 1.695   | 1.920               | 1.924   | 1.972               |        |                     | 1.851   | 6.7   |
| 1,3,5-Trimethylbenzene    | 1.517   | 1.841               | 1.851   | 1.894               |        |                     | 1.735   | 10.2  |
| 1,2,4-Trimethylbenzene    | 1.671   | 2.168               | 2.245   | 2.348               |        |                     | 2.026   | 15.7  |
| 1,3-Dichlorobenzene       | 0.948   | 1.119               | 1.121   | 1.119               |        |                     | 1.053   | 8.6   |
| 1,4-Dichlorobenzene       | 0.944   | 1.122               | 1.115   | 1.138               |        |                     | 1.059   | 8.7   |
| 1,2-Dichlorobenzene       | 0.904   | 1.114               | 1.093   | 1.085               |        |                     | 1.024   | 9.9   |
| 1,2,4-Trichlorobenzene    | 0.905   | 1.224               | 1.117   | 1.026               |        |                     | 1.040   | 12.8  |
| Hexachloro-1,3-Butadiene  | 0.863   | 1.135               | 1.136   | 1.163               |        |                     | 1.034   | 14.7  |
| 1,3-Butadiene             | 0.651   | 0.735               | 0.836   | 1.015               |        |                     | 0.783   | 18.9  |
| Naphthalene               | 2.004   | 2.909               | 2.756   | 2.395               | 2.433  |                     | 2.423   | 15.1  |
| 4-Ethyltoluene            | 1.765   | 2.054               | 2.118   | 2.055               |        |                     | 1.965   | 7.9   |
| 1-Bromo-4-Fluorobenzene   | 0.831   | 0.822               | 0.818   | 0.824               | 0.814  | 0.819               | 0.822   | 0.7   |
| Hexane                    | 1.424   | 1.625               | 1.719   | 1.899               |        |                     | 1.633   | 11.5  |
| Allyl Chloride            | 0.973   | 1.110               | 1.213   | 1.276               |        |                     | 1.130   | 10.4  |
| 1,4-Dioxane               | 181.866 | 232.301             | 242.757 | 262.618             |        |                     | 220.893 | 16.3  |

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 Heated Purge: (Y/N) N Calibration Time(s): 09:44 12:54  
 GC Column: RTX-1 ID: 0.32 (mm)

|                     |        |                     |        |                     |        |                     |       |       |
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|                     |        | RRF0.5 = VL042352.D |        | RRF0.1 = VL042353.D |        | RRF.03 = VL042354.D |       |       |
| COMPOUND            | RRF010 | RRF002              | RRF001 | RRF0.5              | RRF0.1 | RRF.03              | RRF   | % RSD |
| Methyl Methacrylate | 0.430  | 0.495               | 0.520  | 0.545               |        |                     | 0.486 | 10.2  |

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