

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: ROYF02

Lab Code: ACE

SDG No.: Q3321

Instrument ID: MSVOA\_Y

Calibration Date(s): 10/07/2025 10/07/2025

Heated Purge: (Y/N) Y

Calibration Time(s): 09:17 12:32

GC Column: RXI-624 ID: 0.25 (mm)

| LAB FILE ID:                   |        | RRF005 = VY023384.D |        | RRF010 = VY023385.D |        | RRF020 = VY023386.D |       |       |  |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                                |        | RRF050 = VY023387.D |        | RRF100 = VY023388.D |        | RRF150 = VY023389.D |       |       |  |
| COMPOUND                       | RRF005 | RRF010              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |  |
| Dichlorodifluoromethane        | 0.397  | 0.399               | 0.400  | 0.321               | 0.310  | 0.333               | 0.360 | 12    |  |
| Chloromethane                  | 0.518  | 0.557               | 0.512  | 0.449               | 0.438  | 0.467               | 0.490 | 9.4   |  |
| Vinyl Chloride                 | 0.660  | 0.684               | 0.659  | 0.598               | 0.597  | 0.633               | 0.639 | 5.6   |  |
| Bromomethane                   | 0.601  | 0.596               | 0.548  | 0.469               | 0.483  | 0.518               | 0.536 | 10.4  |  |
| Chloroethane                   | 0.441  | 0.462               | 0.453  | 0.414               | 0.413  | 0.429               | 0.435 | 4.7   |  |
| Trichlorofluoromethane         | 0.884  | 0.920               | 0.946  | 0.842               | 0.850  | 0.877               | 0.887 | 4.5   |  |
| 1,1,2-Trichlorotrifluoroethane | 0.484  | 0.486               | 0.479  | 0.437               | 0.429  | 0.445               | 0.460 | 5.6   |  |
| 1,1-Dichloroethene             | 0.449  | 0.475               | 0.456  | 0.427               | 0.426  | 0.443               | 0.446 | 4.1   |  |
| Acetone                        | 0.128  | 0.117               | 0.094  | 0.084               | 0.085  | 0.089               | 0.100 | 18.7  |  |
| Carbon Disulfide               | 1.291  | 1.367               | 1.320  | 1.225               | 1.210  | 1.246               | 1.277 | 4.7   |  |
| Methyl tert-butyl Ether        | 1.006  | 1.075               | 1.075  | 1.020               | 1.091  | 1.165               | 1.072 | 5.3   |  |
| Methyl Acetate                 | 0.244  | 0.265               | 0.260  | 0.245               | 0.266  | 0.286               | 0.261 | 6     |  |
| Methylene Chloride             | 0.625  | 0.634               | 0.576  | 0.458               | 0.463  | 0.473               | 0.538 | 15.4  |  |
| trans-1,2-Dichloroethene       | 0.494  | 0.504               | 0.507  | 0.470               | 0.478  | 0.494               | 0.491 | 2.9   |  |
| 1,1-Dichloroethane             | 0.800  | 0.835               | 0.827  | 0.759               | 0.765  | 0.792               | 0.796 | 3.9   |  |
| Cyclohexane                    | 0.797  | 0.717               | 0.697  | 0.643               | 0.644  | 0.676               | 0.696 | 8.3   |  |
| 2-Butanone                     | 0.133  | 0.134               | 0.115  | 0.105               | 0.115  | 0.126               | 0.122 | 9.6   |  |
| Carbon Tetrachloride           | 0.483  | 0.501               | 0.504  | 0.481               | 0.487  | 0.503               | 0.493 | 2.1   |  |
| cis-1,2-Dichloroethene         | 0.542  | 0.584               | 0.582  | 0.542               | 0.563  | 0.591               | 0.567 | 3.8   |  |
| Bromochloromethane             | 0.332  | 0.323               | 0.320  | 0.267               | 0.270  | 0.273               | 0.297 | 10.2  |  |
| Chloroform                     | 0.886  | 0.912               | 0.901  | 0.835               | 0.842  | 0.874               | 0.875 | 3.5   |  |
| 1,1,1-Trichloroethane          | 0.808  | 0.814               | 0.822  | 0.767               | 0.764  | 0.796               | 0.795 | 3.1   |  |
| Methylcyclohexane              | 0.491  | 0.486               | 0.537  | 0.527               | 0.554  | 0.583               | 0.530 | 7     |  |
| Benzene                        | 1.253  | 1.302               | 1.322  | 1.259               | 1.281  | 1.312               | 1.288 | 2.2   |  |
| 1,2-Dichloroethane             | 0.328  | 0.346               | 0.343  | 0.315               | 0.324  | 0.338               | 0.332 | 3.6   |  |
| Trichloroethene                | 0.380  | 0.381               | 0.386  | 0.367               | 0.372  | 0.376               | 0.377 | 1.8   |  |
| 1,2-Dichloropropane            | 0.276  | 0.289               | 0.292  | 0.274               | 0.283  | 0.293               | 0.285 | 2.9   |  |
| Bromodichloromethane           | 0.440  | 0.463               | 0.459  | 0.439               | 0.449  | 0.468               | 0.453 | 2.7   |  |
| 4-Methyl-2-Pentanone           | 0.141  | 0.159               | 0.161  | 0.159               | 0.180  | 0.196               | 0.166 | 11.6  |  |
| Toluene                        | 0.777  | 0.818               | 0.855  | 0.827               | 0.855  | 0.885               | 0.836 | 4.5   |  |

\* 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.

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                                                |                                                          |
|------------------------------------------------|----------------------------------------------------------|
| Lab Name: <u>Alliance</u>                      | Contract: <u>ROYF02</u>                                  |
| Lab Code: <u>ACE</u>                           | SDG No.: <u>Q3321</u>                                    |
| Instrument ID: <u>MSVOA_Y</u>                  | Calibration Date(s): <u>10/07/2025</u> <u>10/07/2025</u> |
| Heated Purge: (Y/N) <u>Y</u>                   | Calibration Time(s): <u>09:17</u> <u>12:32</u>           |
| GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm) |                                                          |

|                             |        |                     |        |                     |        |                     |       |       |
|-----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                |        | RRF005 = VY023384.D |        | RRF010 = VY023385.D |        | RRF020 = VY023386.D |       |       |
|                             |        | RRF050 = VY023387.D |        | RRF100 = VY023388.D |        | RRF150 = VY023389.D |       |       |
| COMPOUND                    | RRF005 | RRF010              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |
| t-1,3-Dichloropropene       | 0.357  | 0.380               | 0.390  | 0.385               | 0.415  | 0.438               | 0.394 | 7.2   |
| cis-1,3-Dichloropropene     | 0.424  | 0.460               | 0.465  | 0.455               | 0.481  | 0.505               | 0.465 | 5.9   |
| 1,1,2-Trichloroethane       | 0.234  | 0.247               | 0.241  | 0.231               | 0.244  | 0.254               | 0.242 | 3.5   |
| 2-Hexanone                  | 0.108  | 0.117               | 0.110  | 0.112               | 0.126  | 0.138               | 0.119 | 9.9   |
| Dibromochloromethane        | 0.319  | 0.337               | 0.340  | 0.327               | 0.347  | 0.365               | 0.339 | 4.7   |
| 1,2-Dibromoethane           | 0.215  | 0.229               | 0.231  | 0.223               | 0.237  | 0.249               | 0.231 | 5.2   |
| Tetrachloroethene           | 0.538  | 0.517               | 0.561  | 0.523               | 0.502  | 0.456               | 0.516 | 6.9   |
| Chlorobenzene               | 1.086  | 1.095               | 1.092  | 1.053               | 1.083  | 1.116               | 1.087 | 1.9   |
| Ethyl Benzene               | 1.606  | 1.711               | 1.753  | 1.777               | 1.824  | 1.891               | 1.760 | 5.5   |
| m/p-Xylenes                 | 0.632  | 0.675               | 0.715  | 0.719               | 0.734  | 0.763               | 0.706 | 6.5   |
| o-Xylene                    | 0.588  | 0.623               | 0.653  | 0.670               | 0.701  | 0.738               | 0.662 | 8.1   |
| Styrene                     | 0.902  | 1.026               | 1.098  | 1.109               | 1.166  | 1.233               | 1.089 | 10.5  |
| Bromoform                   | 0.225  | 0.242               | 0.238  | 0.230               | 0.244  | 0.262               | 0.240 | 5.4   |
| Isopropylbenzene            | 2.967  | 3.153               | 3.353  | 3.317               | 3.425  | 3.555               | 3.295 | 6.3   |
| 1,1,2,2-Tetrachloroethane   | 0.530  | 0.548               | 0.502  | 0.499               | 0.543  | 0.611               | 0.539 | 7.6   |
| 1,3-Dichlorobenzene         | 1.667  | 1.698               | 1.718  | 1.657               | 1.701  | 1.794               | 1.706 | 2.9   |
| 1,4-Dichlorobenzene         | 1.703  | 1.703               | 1.694  | 1.624               | 1.653  | 1.732               | 1.685 | 2.3   |
| 1,2-Dichlorobenzene         | 1.485  | 1.498               | 1.511  | 1.446               | 1.490  | 1.543               | 1.496 | 2.1   |
| 1,2-Dibromo-3-Chloropropane | 0.087  | 0.093               | 0.083  | 0.083               | 0.088  | 0.096               | 0.088 | 6.2   |
| 1,2,4-Trichlorobenzene      | 0.846  | 0.867               | 0.901  | 0.918               | 1.007  | 1.054               | 0.932 | 8.7   |
| 1,2,3-Trichlorobenzene      | 0.719  | 0.753               | 0.795  | 0.793               | 0.866  | 0.920               | 0.808 | 9.1   |
| 1,2-Dichloroethane-d4       | 0.455  | 0.436               | 0.429  | 0.392               | 0.392  | 0.405               | 0.418 | 6.2   |
| Dibromofluoromethane        | 0.311  | 0.314               | 0.317  | 0.298               | 0.299  | 0.303               | 0.307 | 2.6   |
| Toluene-d8                  | 1.132  | 1.170               | 1.177  | 1.129               | 1.122  | 1.135               | 1.144 | 2     |
| 4-Bromofluorobenzene        | 0.401  | 0.370               | 0.373  | 0.365               | 0.372  | 0.386               | 0.378 | 3.5   |

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