

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: Alliance

Contract: CHEM02

Lab Code: ACE

SDG No.: Q3745

Instrument ID: MSVOA\_N

Calibration Date(s): 11/24/2025 11/24/2025

Heated Purge: (Y/N) N

Calibration Time(s): 12:38 14:35

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

| LAB FILE ID:                   |        | RRF001 = VN088344.D |        | RRF005 = VN088345.D |        | RRF020 = VN088346.D |       |       |  |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                                |        | RRF050 = VN088347.D |        | RRF100 = VN088348.D |        | RRF150 = VN088349.D |       |       |  |
| COMPOUND                       | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |  |
| Dichlorodifluoromethane        | 0.719  | 0.676               | 0.796  | 0.741               | 0.838  | 0.736               | 0.751 | 7.7   |  |
| Chloromethane                  | 0.911  | 0.800               | 0.805  | 0.801               | 0.759  | 0.764               | 0.807 | 6.8   |  |
| Vinyl Chloride                 | 0.804  | 0.844               | 0.856  | 0.812               | 0.837  | 0.772               | 0.821 | 3.8   |  |
| Ethyl Acetate                  | 0.735  | 0.678               | 0.724  | 0.720               | 0.681  | 0.681               | 0.703 | 3.7   |  |
| Bromomethane                   |        | 0.407               | 0.392  | 0.396               | 0.387  | 0.392               | 0.395 | 1.9   |  |
| Chloroethane                   | 0.555  | 0.503               | 0.496  | 0.505               | 0.502  | 0.485               | 0.508 | 4.8   |  |
| Trichlorofluoromethane         | 1.234  | 1.170               | 1.168  | 1.092               | 1.201  | 1.055               | 1.153 | 5.8   |  |
| 1,1,2-Trichlorotrifluoroethane | 0.660  | 0.635               | 0.639  | 0.564               | 0.633  | 0.553               | 0.614 | 7.2   |  |
| Tert butyl alcohol             |        | 0.151               | 0.149  | 0.161               | 0.142  | 0.148               | 0.150 | 4.6   |  |
| 1,1-Dichloroethene             | 0.722  | 0.620               | 0.601  | 0.568               | 0.595  | 0.554               | 0.610 | 9.8   |  |
| Acrolein                       |        | 0.146               | 0.132  | 0.151               | 0.143  | 0.147               | 0.144 | 5     |  |
| Acrylonitrile                  | 0.435  | 0.435               | 0.432  | 0.466               | 0.416  | 0.418               | 0.434 | 4.1   |  |
| Acetone                        | 0.412  | 0.355               | 0.322  | 0.335               | 0.305  | 0.314               | 0.341 | 11.5  |  |
| Carbon Disulfide               | 2.061  | 2.043               | 2.043  | 1.983               | 1.996  | 1.833               | 1.993 | 4.2   |  |
| Methyl tert-butyl Ether        | 2.126  | 2.239               | 2.245  | 2.504               | 2.265  | 2.333               | 2.285 | 5.5   |  |
| Methyl Acetate                 | 1.072  | 1.048               | 0.957  | 1.049               | 0.972  | 0.985               | 1.014 | 4.7   |  |
| Methylene Chloride             | 0.797  | 0.752               | 0.704  | 0.756               | 0.686  | 0.678               | 0.729 | 6.4   |  |
| trans-1,2-Dichloroethene       | 0.752  | 0.668               | 0.676  | 0.686               | 0.647  | 0.626               | 0.676 | 6.4   |  |
| Vinyl Acetate                  | 1.711  | 1.879               | 1.997  | 2.137               | 1.943  | 2.002               | 1.945 | 7.4   |  |
| 1,1-Dichloroethane             | 1.389  | 1.431               | 1.348  | 1.427               | 1.321  | 1.286               | 1.367 | 4.3   |  |
| Cyclohexane                    |        | 1.450               | 1.264  | 1.133               | 1.237  | 1.109               | 1.239 | 10.9  |  |
| 2-Butanone                     | 0.509  | 0.538               | 0.548  | 0.596               | 0.540  | 0.556               | 0.548 | 5.2   |  |
| Carbon Tetrachloride           | 0.565  | 0.504               | 0.548  | 0.499               | 0.570  | 0.509               | 0.532 | 6     |  |
| 2,2-Dichloropropane            | 1.245  | 1.133               | 1.185  | 1.204               | 1.176  | 1.137               | 1.180 | 3.6   |  |
| cis-1,2-Dichloroethene         | 0.797  | 0.801               | 0.794  | 0.829               | 0.747  | 0.744               | 0.785 | 4.2   |  |
| Bromochloromethane             | 0.724  | 0.722               | 0.684  | 0.735               | 0.624  | 0.597               | 0.681 | 8.5   |  |
| Chloroform                     | 1.377  | 1.362               | 1.351  | 1.392               | 1.291  | 1.285               | 1.343 | 3.3   |  |
| 1,1,1-Trichloroethane          | 1.247  | 1.219               | 1.192  | 1.158               | 1.166  | 1.104               | 1.181 | 4.3   |  |
| Methylcyclohexane              | 0.534  | 0.542               | 0.606  | 0.527               | 0.662  | 0.561               | 0.572 | 9.1   |  |
| 1,1-Dichloropropene            | 0.607  | 0.542               | 0.531  | 0.491               | 0.547  | 0.490               | 0.535 | 8.1   |  |

\* 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>CHEM02</u>                                  |
| Lab Code: <u>ACE</u>                           | SDG No.: <u>Q3745</u>                                    |
| Instrument ID: <u>MSVOA_N</u>                  | Calibration Date(s): <u>11/24/2025</u> <u>11/24/2025</u> |
| Heated Purge: (Y/N) <u>N</u>                   | Calibration Time(s): <u>12:38</u> <u>14:35</u>           |
| GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm) |                                                          |

|                           |                     |        |        |                     |        |                     |       |       |
|---------------------------|---------------------|--------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:              | RRF001 = VN088344.D |        |        | RRF005 = VN088345.D |        | RRF020 = VN088346.D |       |       |
|                           | RRF050 = VN088347.D |        |        | RRF100 = VN088348.D |        | RRF150 = VN088349.D |       |       |
| COMPOUND                  | RRF001              | RRF005 | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |
| Benzene                   | 1.658               | 1.555  | 1.609  | 1.559               | 1.559  | 1.477               | 1.570 | 3.9   |
| 1,2-Dichloroethane        | 0.583               | 0.614  | 0.596  | 0.615               | 0.597  | 0.587               | 0.599 | 2.3   |
| Trichloroethene           | 0.382               | 0.385  | 0.383  | 0.357               | 0.371  | 0.348               | 0.371 | 4.1   |
| 1,2-Dichloropropane       | 0.426               | 0.402  | 0.402  | 0.406               | 0.396  | 0.381               | 0.402 | 3.6   |
| Dibromomethane            | 0.288               | 0.292  | 0.289  | 0.293               | 0.283  | 0.277               | 0.287 | 2.1   |
| Bromodichloromethane      | 0.602               | 0.586  | 0.586  | 0.589               | 0.583  | 0.566               | 0.586 | 2     |
| 4-Methyl-2-Pentanone      | 0.540               | 0.616  | 0.661  | 0.673               | 0.658  | 0.651               | 0.633 | 7.8   |
| Toluene                   | 0.829               | 0.885  | 0.940  | 0.913               | 0.960  | 0.911               | 0.906 | 5.1   |
| t-1,3-Dichloropropene     | 0.567               | 0.553  | 0.598  | 0.631               | 0.631  | 0.625               | 0.601 | 5.7   |
| cis-1,3-Dichloropropene   | 0.590               | 0.618  | 0.628  | 0.662               | 0.654  | 0.640               | 0.632 | 4.2   |
| 1,1,2-Trichloroethane     | 0.350               | 0.359  | 0.364  | 0.353               | 0.341  | 0.336               | 0.351 | 3     |
| 1,3-Dichloropropane       | 0.607               | 0.638  | 0.647  | 0.644               | 0.638  | 0.622               | 0.633 | 2.4   |
| 2-Chloroethyl Vinyl ether | 0.258               | 0.323  | 0.333  | 0.364               | 0.337  | 0.301               | 0.319 | 11.4  |
| 2-Hexanone                | 0.359               | 0.301  | 0.406  | 0.436               | 0.442  | 0.450               | 0.399 | 14.7  |
| Dibromochloromethane      | 0.384               | 0.386  | 0.406  | 0.409               | 0.410  | 0.404               | 0.400 | 2.9   |
| 1,2-Dibromoethane         | 0.304               | 0.367  | 0.364  | 0.369               | 0.367  | 0.359               | 0.355 | 7.2   |
| Tetrachloroethene         | 0.426               | 0.388  | 0.413  | 0.377               | 0.407  | 0.367               | 0.396 | 5.7   |
| Chlorobenzene             | 1.114               | 1.142  | 1.161  | 1.106               | 1.145  | 1.070               | 1.123 | 3     |
| 1,1,1,2-Tetrachloroethane | 0.378               | 0.372  | 0.394  | 0.383               | 0.386  | 0.361               | 0.379 | 3     |
| Hexachloroethane          | 0.549               | 0.550  | 0.571  | 0.535               | 0.604  | 0.553               | 0.560 | 4.4   |
| Ethyl Benzene             | 1.881               | 1.860  | 2.014  | 1.938               | 2.133  | 1.952               | 1.963 | 5.1   |
| m/p-Xylenes               | 0.658               | 0.690  | 0.769  | 0.733               | 0.788  | 0.728               | 0.728 | 6.6   |
| o-Xylene                  | 0.670               | 0.619  | 0.722  | 0.696               | 0.749  | 0.708               | 0.694 | 6.5   |
| Styrene                   | 0.884               | 1.025  | 1.163  | 1.193               | 1.284  | 1.210               | 1.127 | 12.9  |
| Bromoform                 | 0.224               | 0.273  | 0.285  | 0.285               | 0.302  | 0.295               | 0.277 | 10.1  |
| Isopropylbenzene          | 3.339               | 3.528  | 3.944  | 3.624               | 4.040  | 3.706               | 3.697 | 7.1   |
| 1,1,2,2-Tetrachloroethane | 1.377               | 1.218  | 1.231  | 1.127               | 1.135  | 1.082               | 1.195 | 8.9   |
| 1,2,3-Trichloropropane    | 1.036               | 1.199  | 1.268  | 1.202               | 1.177  | 1.145               | 1.171 | 6.6   |
| Bromobenzene              | 0.781               | 0.885  | 0.866  | 0.863               | 0.865  | 0.816               | 0.846 | 4.7   |
| n-propylbenzene           | 4.076               | 4.270  | 4.925  | 4.439               | 4.931  | 4.496               | 4.523 | 7.7   |

\* 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.

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                                                |                                                          |
|------------------------------------------------|----------------------------------------------------------|
| Lab Name: <u>Alliance</u>                      | Contract: <u>CHEM02</u>                                  |
| Lab Code: <u>ACE</u>                           | SDG No.: <u>Q3745</u>                                    |
| Instrument ID: <u>MSVOA_N</u>                  | Calibration Date(s): <u>11/24/2025</u> <u>11/24/2025</u> |
| Heated Purge: (Y/N) <u>N</u>                   | Calibration Time(s): <u>12:38</u> <u>14:35</u>           |
| GC Column: <u>RXI-624</u> ID: <u>0.25</u> (mm) |                                                          |

|                             |        |                     |        |                     |        |                     |       |       |
|-----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                |        | RRF001 = VN088344.D |        | RRF005 = VN088345.D |        | RRF020 = VN088346.D |       |       |
|                             |        | RRF050 = VN088347.D |        | RRF100 = VN088348.D |        | RRF150 = VN088349.D |       |       |
| COMPOUND                    | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |
| 2-Chlorotoluene             | 2.707  | 2.799               | 2.851  | 2.699               | 2.835  | 2.681               | 2.762 | 2.7   |
| 1,3,5-Trimethylbenzene      | 2.650  | 2.988               | 3.315  | 3.051               | 3.307  | 3.069               | 3.063 | 8     |
| 4-Chlorotoluene             | 2.093  | 2.830               | 2.935  | 2.854               | 2.944  | 2.803               | 2.743 | 11.8  |
| tert-Butylbenzene           | 2.509  | 2.468               | 2.773  | 2.509               | 2.816  | 2.549               | 2.604 | 5.8   |
| 1,2,4-Trimethylbenzene      | 2.572  | 2.827               | 3.310  | 3.112               | 3.344  | 3.105               | 3.045 | 9.7   |
| sec-Butylbenzene            | 3.457  | 3.614               | 3.971  | 3.609               | 4.122  | 3.674               | 3.741 | 6.7   |
| p-Isopropyltoluene          | 2.714  | 2.808               | 3.330  | 3.047               | 3.435  | 3.073               | 3.068 | 9.2   |
| 1,3-Dichlorobenzene         | 1.730  | 1.573               | 1.720  | 1.606               | 1.676  | 1.574               | 1.647 | 4.3   |
| 1,4-Dichlorobenzene         | 1.747  | 1.825               | 1.786  | 1.681               | 1.702  | 1.608               | 1.725 | 4.5   |
| n-Butylbenzene              | 2.827  | 2.664               | 3.155  | 2.873               | 3.373  | 2.979               | 2.979 | 8.5   |
| 1,2-Dichlorobenzene         | 1.518  | 1.574               | 1.583  | 1.521               | 1.595  | 1.505               | 1.550 | 2.5   |
| 1,2-Dibromo-3-Chloropropane | 0.284  | 0.284               | 0.284  | 0.288               | 0.300  | 0.300               | 0.290 | 2.8   |
| 1,2,4-Trichlorobenzene      | 0.998  | 0.864               | 0.901  | 0.889               | 0.955  | 0.917               | 0.921 | 5.3   |
| Hexachlorobutadiene         | 0.362  | 0.316               | 0.349  | 0.297               | 0.354  | 0.307               | 0.331 | 8.3   |
| Naphthalene                 | 2.946  | 2.864               | 3.163  | 3.292               | 3.483  | 3.395               | 3.190 | 7.7   |
| 1,2,3-Trichlorobenzene      | 0.942  | 0.850               | 0.888  | 0.875               | 0.930  | 0.891               | 0.896 | 3.8   |
| 1,2-Dichloroethane-d4       |        | 0.954               | 0.916  | 0.999               | 0.882  | 0.960               | 0.942 | 4.7   |
| Dibromofluoromethane        |        | 0.361               | 0.350  | 0.350               | 0.334  | 0.337               | 0.346 | 3.2   |
| Toluene-d8                  |        | 1.249               | 1.292  | 1.285               | 1.308  | 1.313               | 1.289 | 2     |
| 4-Bromofluorobenzene        |        | 0.402               | 0.413  | 0.448               | 0.466  | 0.482               | 0.442 | 7.7   |

\* 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.