

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

|                     |                 |                      |                                     |
|---------------------|-----------------|----------------------|-------------------------------------|
| Lab Name:           | <u>Alliance</u> | Contract:            | <u>AECO15</u>                       |
| Lab Code:           | <u>ACE</u>      | SDG No.:             | <u>Q3193</u>                        |
| Instrument ID:      | <u>MSVOA_X</u>  | Calibration Date(s): | <u>09/16/2025</u> <u>09/16/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>        | Calibration Time(s): | <u>09:23</u> <u>12:01</u>           |
| GC Column:          | <u>DB-624UI</u> | ID:                  | <u>0.18</u> (mm)                    |

| LAB FILE ID:                   |        | RRF001 = VX047585.D |        | RRF005 = VX047586.D |        | RRF020 = VX047587.D |       |       |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
|                                |        | RRF050 = VX047588.D |        | RRF100 = VX047589.D |        | RRF150 = VX047590.D |       |       |
| COMPOUND                       | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |
| Dichlorodifluoromethane        | 0.447  | 0.575               | 0.582  | 0.501               | 0.510  | 0.492               | 0.518 | 10    |
| Chloromethane                  | 0.610  | 0.749               | 0.739  | 0.661               | 0.677  | 0.647               | 0.680 | 7.9   |
| Vinyl Chloride                 | 0.572  | 0.713               | 0.716  | 0.661               | 0.677  | 0.658               | 0.666 | 7.9   |
| Bromomethane                   |        | 0.493               | 0.451  | 0.424               | 0.423  | 0.321               | 0.422 | 15    |
| Chloroethane                   | 0.433  | 0.450               | 0.469  | 0.443               | 0.450  | 0.426               | 0.445 | 3.4   |
| Trichlorofluoromethane         | 0.865  | 1.033               | 1.038  | 0.995               | 1.018  | 0.986               | 0.989 | 6.5   |
| 1,1,2-Trichlorotrifluoroethane | 0.519  | 0.591               | 0.603  | 0.583               | 0.590  | 0.583               | 0.578 | 5.2   |
| 1,1-Dichloroethene             | 0.526  | 0.612               | 0.618  | 0.604               | 0.609  | 0.596               | 0.594 | 5.7   |
| Acetone                        | 0.343  | 0.416               | 0.400  | 0.317               | 0.302  | 0.299               | 0.346 | 14.7  |
| Carbon Disulfide               | 1.555  | 1.735               | 1.758  | 1.566               | 1.589  | 1.555               | 1.626 | 5.8   |
| Methyl tert-butyl Ether        | 1.884  | 2.238               | 2.221  | 2.247               | 2.212  | 2.209               | 2.169 | 6.5   |
| Methyl Acetate                 | 0.612  | 0.673               | 0.714  | 0.876               | 0.865  | 0.881               | 0.770 | 15.4  |
| Methylene Chloride             | 0.718  | 0.772               | 0.754  | 0.730               | 0.716  | 0.705               | 0.732 | 3.5   |
| trans-1,2-Dichloroethene       | 0.543  | 0.703               | 0.665  | 0.649               | 0.643  | 0.635               | 0.640 | 8.3   |
| 1,1-Dichloroethane             | 1.090  | 1.325               | 1.299  | 1.297               | 1.287  | 1.278               | 1.263 | 6.8   |
| Cyclohexane                    |        | 1.161               | 1.092  | 1.030               | 0.984  | 0.993               | 1.052 | 7.1   |
| 2-Butanone                     | 0.370  | 0.470               | 0.473  | 0.439               | 0.414  | 0.424               | 0.432 | 9     |
| Carbon Tetrachloride           | 0.482  | 0.574               | 0.555  | 0.539               | 0.526  | 0.518               | 0.532 | 6     |
| cis-1,2-Dichloroethene         | 0.698  | 0.806               | 0.803  | 0.808               | 0.787  | 0.797               | 0.783 | 5.4   |
| Bromochloromethane             | 0.457  | 0.607               | 0.450  | 0.577               | 0.587  | 0.630               | 0.551 | 14.2  |
| Chloroform                     | 1.034  | 1.364               | 1.366  | 1.323               | 1.282  | 1.288               | 1.276 | 9.7   |
| 1,1,1-Trichloroethane          | 0.901  | 1.123               | 1.141  | 1.127               | 1.100  | 1.103               | 1.082 | 8.3   |
| Methylcyclohexane              | 0.490  | 0.589               | 0.591  | 0.565               | 0.548  | 0.554               | 0.556 | 6.6   |
| Benzene                        | 1.330  | 1.577               | 1.566  | 1.523               | 1.473  | 1.459               | 1.488 | 6.1   |
| 1,2-Dichloroethane             | 0.513  | 0.595               | 0.600  | 0.582               | 0.554  | 0.553               | 0.566 | 5.8   |
| Trichloroethene                | 0.296  | 0.382               | 0.382  | 0.370               | 0.364  | 0.363               | 0.360 | 8.9   |
| 1,2-Dichloropropane            | 0.326  | 0.403               | 0.394  | 0.397               | 0.382  | 0.384               | 0.381 | 7.4   |
| Bromodichloromethane           | 0.461  | 0.599               | 0.601  | 0.607               | 0.579  | 0.586               | 0.572 | 9.7   |
| 4-Methyl-2-Pentanone           | 0.390  | 0.493               | 0.503  | 0.511               | 0.479  | 0.487               | 0.477 | 9.3   |
| Toluene                        | 0.841  | 0.967               | 0.960  | 0.940               | 0.897  | 0.897               | 0.917 | 5.2   |

\* 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>AECO15</u>                       |
| Lab Code:           | <u>ACE</u>                           | SDG No.:             | <u>Q3193</u>                        |
| Instrument ID:      | <u>MSVOA_X</u>                       | Calibration Date(s): | <u>09/16/2025</u> <u>09/16/2025</u> |
| Heated Purge: (Y/N) | <u>N</u>                             | Calibration Time(s): | <u>09:23</u> <u>12:01</u>           |
| GC Column:          | <u>DB-624UI</u> ID: <u>0.18</u> (mm) |                      |                                     |

|                             |        |                     |        |                     |        |                     |       |       |
|-----------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|
| LAB FILE ID:                |        | RRF001 = VX047585.D |        | RRF005 = VX047586.D |        | RRF020 = VX047587.D |       |       |
|                             |        | RRF050 = VX047588.D |        | RRF100 = VX047589.D |        | RRF150 = VX047590.D |       |       |
| COMPOUND                    | RRF001 | RRF005              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |
| t-1,3-Dichloropropene       | 0.487  | 0.588               | 0.602  | 0.619               | 0.595  | 0.611               | 0.584 | 8.3   |
| cis-1,3-Dichloropropene     | 0.513  | 0.652               | 0.644  | 0.653               | 0.637  | 0.644               | 0.624 | 8.7   |
| 1,1,2-Trichloroethane       | 0.323  | 0.365               | 0.371  | 0.367               | 0.344  | 0.351               | 0.354 | 5.1   |
| 2-Hexanone                  | 0.274  | 0.374               | 0.372  | 0.360               | 0.333  | 0.343               | 0.343 | 10.8  |
| Dibromochloromethane        | 0.321  | 0.414               | 0.431  | 0.435               | 0.421  | 0.422               | 0.407 | 10.5  |
| 1,2-Dibromoethane           | 0.287  | 0.379               | 0.384  | 0.379               | 0.364  | 0.367               | 0.360 | 10.1  |
| Tetrachloroethene           | 0.299  | 0.354               | 0.342  | 0.326               | 0.320  | 0.314               | 0.326 | 6     |
| Chlorobenzene               | 0.998  | 1.182               | 1.177  | 1.178               | 1.139  | 1.142               | 1.136 | 6.2   |
| Ethyl Benzene               | 1.625  | 2.077               | 2.049  | 2.039               | 1.996  | 1.973               | 1.960 | 8.6   |
| m/p-Xylenes                 | 0.604  | 0.774               | 0.765  | 0.765               | 0.739  | 0.730               | 0.729 | 8.8   |
| o-Xylene                    | 0.576  | 0.734               | 0.734  | 0.749               | 0.721  | 0.719               | 0.706 | 9.1   |
| Styrene                     | 1.019  | 1.288               | 1.290  | 1.299               | 1.257  | 1.265               | 1.236 | 8.7   |
| Bromoform                   | 0.221  | 0.293               | 0.301  | 0.313               | 0.312  | 0.317               | 0.293 | 12.4  |
| Isopropylbenzene            | 3.124  | 3.907               | 3.986  | 4.047               | 3.907  | 3.839               | 3.801 | 8.9   |
| 1,1,2,2-Tetrachloroethane   | 0.965  | 1.182               | 1.203  | 1.225               | 1.152  | 1.173               | 1.150 | 8.2   |
| 1,3-Dichlorobenzene         | 1.462  | 1.800               | 1.829  | 1.817               | 1.768  | 1.758               | 1.739 | 8     |
| 1,4-Dichlorobenzene         | 1.570  | 1.890               | 1.868  | 1.832               | 1.767  | 1.783               | 1.785 | 6.5   |
| 1,2-Dichlorobenzene         | 1.359  | 1.691               | 1.757  | 1.746               | 1.696  | 1.692               | 1.657 | 9     |
| 1,2-Dibromo-3-Chloropropane | 0.173  | 0.218               | 0.242  | 0.254               | 0.250  | 0.260               | 0.233 | 14.1  |
| 1,2,4-Trichlorobenzene      | 0.831  | 1.099               | 1.107  | 1.156               | 1.140  | 1.126               | 1.077 | 11.3  |
| 1,2,3-Trichlorobenzene      | 0.753  | 0.986               | 1.031  | 1.088               | 1.078  | 1.073               | 1.002 | 12.7  |
| 1,2-Dichloroethane-d4       |        | 0.884               | 0.647  | 0.770               | 0.815  | 0.863               | 0.796 | 11.8  |
| Dibromofluoromethane        |        | 0.356               | 0.266  | 0.331               | 0.355  | 0.368               | 0.335 | 12.3  |
| Toluene-d8                  |        | 1.237               | 0.943  | 1.147               | 1.211  | 1.263               | 1.160 | 11.1  |
| 4-Bromofluorobenzene        |        | 0.478               | 0.360  | 0.427               | 0.452  | 0.484               | 0.440 | 11.4  |

\* 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.