

### VOLATILE ORGANICS INITIAL CALIBRATION DATA

Lab Name: CHEMTECH Contract: ROMA02

Lab Code: CHEM Case No.: P4258 SAS No.: P4258 SDG No.: P4258

Instrument ID: MSVOA\_Y Calibration Date(s): 09/09/2024 09/09/2024

Heated Purge: (Y/N) Y Calibration Time(s): 09:52 11:53

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

| LAB FILE ID:                   |        | RRF005 = VY019454.D |        | RRF010 = VY019455.D |        | RRF020 = VY019456.D |       |       |  |
|--------------------------------|--------|---------------------|--------|---------------------|--------|---------------------|-------|-------|--|
|                                |        | RRF050 = VY019457.D |        | RRF100 = VY019458.D |        | RRF150 = VY019459.D |       |       |  |
| COMPOUND                       | RRF005 | RRF010              | RRF020 | RRF050              | RRF100 | RRF150              | RRF   | % RSD |  |
| Dichlorodifluoromethane        | 0.375  | 0.371               | 0.373  | 0.265               | 0.248  | 0.264               | 0.316 | 19.8  |  |
| Chloromethane                  | 0.456  | 0.454               | 0.442  | 0.392               | 0.357  | 0.380               | 0.414 | 10.3  |  |
| Vinyl Chloride                 | 0.451  | 0.466               | 0.476  | 0.447               | 0.409  | 0.435               | 0.447 | 5.2   |  |
| Bromomethane                   | 0.255  | 0.278               | 0.273  | 0.283               | 0.269  | 0.291               | 0.275 | 4.6   |  |
| Chloroethane                   | 0.300  | 0.303               | 0.299  | 0.297               | 0.270  | 0.286               | 0.292 | 4.3   |  |
| Trichlorofluoromethane         | 0.760  | 0.775               | 0.776  | 0.733               | 0.686  | 0.743               | 0.746 | 4.6   |  |
| 1,1,2-Trichlorotrifluoroethane | 0.477  | 0.480               | 0.476  | 0.457               | 0.425  | 0.453               | 0.461 | 4.6   |  |
| 1,1-Dichloroethene             | 0.457  | 0.449               | 0.450  | 0.447               | 0.413  | 0.446               | 0.444 | 3.5   |  |
| Acetone                        | 0.102  | 0.113               | 0.106  | 0.124               | 0.107  | 0.109               | 0.110 | 6.9   |  |
| Carbon Disulfide               | 1.144  | 1.159               | 1.161  | 1.283               | 1.177  | 1.266               | 1.199 | 5     |  |
| Methyl tert-butyl Ether        | 1.146  | 1.304               | 1.294  | 1.313               | 1.212  | 1.270               | 1.256 | 5.2   |  |
| Methyl Acetate                 | 0.285  | 0.314               | 0.279  | 0.345               | 0.283  | 0.290               | 0.299 | 8.5   |  |
| Methylene Chloride             | 0.538  | 0.524               | 0.513  | 0.502               | 0.456  | 0.485               | 0.503 | 5.8   |  |
| trans-1,2-Dichloroethene       | 0.488  | 0.488               | 0.484  | 0.500               | 0.461  | 0.496               | 0.486 | 2.9   |  |
| 1,1-Dichloroethane             | 0.889  | 0.945               | 0.936  | 0.945               | 0.862  | 0.923               | 0.917 | 3.7   |  |
| Cyclohexane                    | 0.969  | 0.873               | 0.819  | 0.803               | 0.735  | 0.772               | 0.829 | 10    |  |
| 2-Butanone                     | 0.134  | 0.166               | 0.159  | 0.171               | 0.154  | 0.158               | 0.157 | 8.1   |  |
| Carbon Tetrachloride           | 0.440  | 0.451               | 0.441  | 0.437               | 0.408  | 0.433               | 0.435 | 3.4   |  |
| cis-1,2-Dichloroethene         | 0.568  | 0.592               | 0.602  | 0.613               | 0.564  | 0.605               | 0.591 | 3.4   |  |
| Bromochloromethane             | 0.368  | 0.373               | 0.378  | 0.410               | 0.383  | 0.402               | 0.386 | 4.4   |  |
| Chloroform                     | 0.909  | 0.952               | 0.954  | 0.967               | 0.885  | 0.947               | 0.936 | 3.4   |  |
| 1,1,1-Trichloroethane          | 0.825  | 0.868               | 0.851  | 0.853               | 0.789  | 0.847               | 0.839 | 3.3   |  |
| Methylcyclohexane              | 0.524  | 0.533               | 0.514  | 0.516               | 0.481  | 0.513               | 0.513 | 3.4   |  |
| Benzene                        | 1.169  | 1.235               | 1.211  | 1.212               | 1.118  | 1.184               | 1.188 | 3.5   |  |
| 1,2-Dichloroethane             | 0.297  | 0.336               | 0.325  | 0.330               | 0.307  | 0.324               | 0.320 | 4.7   |  |
| Trichloroethene                | 0.301  | 0.309               | 0.315  | 0.312               | 0.288  | 0.307               | 0.306 | 3.1   |  |
| 1,2-Dichloropropane            | 0.287  | 0.301               | 0.299  | 0.294               | 0.273  | 0.289               | 0.290 | 3.5   |  |
| Bromodichloromethane           | 0.413  | 0.445               | 0.431  | 0.437               | 0.408  | 0.428               | 0.427 | 3.3   |  |
| 4-Methyl-2-Pentanone           | 0.181  | 0.225               | 0.214  | 0.222               | 0.203  | 0.209               | 0.209 | 7.7   |  |
| Toluene                        | 0.738  | 0.793               | 0.775  | 0.790               | 0.725  | 0.764               | 0.764 | 3.6   |  |

\* 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>CHEMTECH</u>  | Contract:            | <u>ROMA02</u>     |
| Lab Code:           | <u>CHEM</u>      | Case No.:            | <u>P4258</u>      |
| Instrument ID:      | <u>MSVOA_Y</u>   | SAS No.:             | <u>P4258</u>      |
| Heated Purge: (Y/N) | <u>Y</u>         | SDG No.:             | <u>P4258</u>      |
| GC Column:          | <u>RXI-624</u>   | Calibration Date(s): | <u>09/09/2024</u> |
| ID:                 | <u>0.25 (mm)</u> | Calibration Time(s): | <u>09:52</u>      |
|                     |                  |                      | <u>11:53</u>      |

| LAB FILE ID:                | RRF005 = VY019454.D | RRF010 = VY019455.D | RRF020 = VY019456.D | RRF050 = VY019457.D | RRF100 = VY019458.D | RRF150 = VY019459.D |       |       |
|-----------------------------|---------------------|---------------------|---------------------|---------------------|---------------------|---------------------|-------|-------|
| COMPOUND                    | RRF005              | RRF010              | RRF020              | RRF050              | RRF100              | RRF150              | RRF   | % RSD |
| t-1,3-Dichloropropene       | 0.366               | 0.424               | 0.416               | 0.426               | 0.393               | 0.417               | 0.407 | 5.8   |
| cis-1,3-Dichloropropene     | 0.444               | 0.491               | 0.485               | 0.492               | 0.453               | 0.478               | 0.474 | 4.3   |
| 1,1,2-Trichloroethane       | 0.209               | 0.227               | 0.219               | 0.231               | 0.211               | 0.220               | 0.219 | 4     |
| 2-Hexanone                  | 0.131               | 0.166               | 0.158               | 0.163               | 0.149               | 0.152               | 0.153 | 8.2   |
| Dibromochloromethane        | 0.260               | 0.302               | 0.297               | 0.304               | 0.281               | 0.296               | 0.290 | 5.8   |
| 1,2-Dibromoethane           | 0.188               | 0.216               | 0.211               | 0.212               | 0.195               | 0.204               | 0.204 | 5.5   |
| Tetrachloroethene           | 0.315               | 0.319               | 0.313               | 0.312               | 0.291               | 0.307               | 0.310 | 3.2   |
| Chlorobenzene               | 0.961               | 1.026               | 1.010               | 0.995               | 0.925               | 0.983               | 0.984 | 3.7   |
| Ethyl Benzene               | 1.739               | 1.856               | 1.796               | 1.769               | 1.626               | 1.701               | 1.748 | 4.5   |
| m/p-Xylenes                 | 0.656               | 0.688               | 0.682               | 0.670               | 0.620               | 0.650               | 0.661 | 3.7   |
| o-Xylene                    | 0.646               | 0.673               | 0.667               | 0.656               | 0.601               | 0.635               | 0.646 | 4     |
| Styrene                     | 1.061               | 1.143               | 1.126               | 1.112               | 1.031               | 1.073               | 1.091 | 3.9   |
| Bromoform                   | 0.173               | 0.207               | 0.196               | 0.205               | 0.189               | 0.198               | 0.195 | 6.3   |
| Isopropylbenzene            | 3.732               | 3.822               | 3.698               | 3.613               | 3.336               | 3.563               | 3.627 | 4.7   |
| 1,1,2,2-Tetrachloroethane   | 0.589               | 0.684               | 0.652               | 0.665               | 0.617               | 0.654               | 0.644 | 5.4   |
| 1,3-Dichlorobenzene         | 1.564               | 1.622               | 1.612               | 1.575               | 1.453               | 1.554               | 1.563 | 3.8   |
| 1,4-Dichlorobenzene         | 1.554               | 1.596               | 1.594               | 1.565               | 1.452               | 1.548               | 1.551 | 3.4   |
| 1,2-Dichlorobenzene         | 1.367               | 1.448               | 1.414               | 1.422               | 1.315               | 1.404               | 1.395 | 3.4   |
| 1,2-Dibromo-3-Chloropropane | 0.099               | 0.116               | 0.105               | 0.110               | 0.103               | 0.110               | 0.107 | 5.5   |
| 1,2,4-Trichlorobenzene      | 0.810               | 0.837               | 0.858               | 0.874               | 0.817               | 0.894               | 0.848 | 3.9   |
| 1,2,3-Trichlorobenzene      | 0.664               | 0.719               | 0.738               | 0.754               | 0.705               | 0.776               | 0.726 | 5.4   |
| 1,2-Dichloroethane-d4       | 0.420               | 0.462               | 0.466               | 0.485               | 0.465               | 0.480               | 0.463 | 5     |
| Dibromofluoromethane        | 0.264               | 0.291               | 0.287               | 0.290               | 0.279               | 0.288               | 0.283 | 3.7   |
| Toluene-d8                  | 0.986               | 1.073               | 1.055               | 1.084               | 1.033               | 1.060               | 1.049 | 3.4   |
| 4-Bromofluorobenzene        | 0.425               | 0.414               | 0.407               | 0.400               | 0.385               | 0.389               | 0.403 | 3.8   |

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