

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q1774 SAS No.: Q1774 SDG No.: Q1774

Instrument ID: MSVOA\_N Calibration Date/Time: 04/11/2025 17:28

Lab File ID: VN086255.D Init. Calib. Date(s): 04/11/2025 04/11/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:04 12:40

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

| COMPOUND                       | RRF   | RRF020 | MIN<br>RRF | %D     | MAX%D |
|--------------------------------|-------|--------|------------|--------|-------|
| Dichlorodifluoromethane        | 2.404 | 2.329  |            | -3.12  | 50    |
| Chloromethane                  | 3.322 | 3.290  |            | -0.96  | 50    |
| Vinyl Chloride                 | 3.164 | 3.093  | 0.1        | -2.24  | 50    |
| Bromomethane                   | 1.578 | 1.524  | 0.1        | -3.42  | 50    |
| Chloroethane                   | 1.998 | 1.989  |            | -0.45  | 50    |
| Trichlorofluoromethane         | 3.433 | 3.320  |            | -3.29  | 50    |
| 1,1,2-Trichlorotrifluoroethane | 2.117 | 1.981  |            | -6.42  | 50    |
| 1,1-Dichloroethene             | 2.284 | 2.142  | 0.1        | -6.22  | 50    |
| Acrolein                       | 0.367 | 0.303  |            | -17.44 | 50    |
| Acrylonitrile                  | 1.362 | 1.283  |            | -5.8   | 50    |
| Acetone                        | 0.345 | 0.315  |            | -8.7   | 50    |
| Carbon Disulfide               | 6.550 | 6.271  |            | -4.26  | 50    |
| Methyl tert-Butyl Ether        | 8.244 | 7.740  |            | -6.11  | 50    |
| Methyl Acetate                 | 3.226 | 2.973  |            | -7.84  | 50    |
| Methylene Chloride             | 2.605 | 2.476  |            | -4.95  | 50    |
| trans-1,2-Dichloroethene       | 2.406 | 2.267  |            | -5.78  | 50    |
| 1,1-Dichloroethane             | 4.567 | 4.406  | 0.2        | -3.53  | 50    |
| Cyclohexane                    | 0.705 | 0.688  |            | -2.41  | 50    |
| 2-Butanone                     | 0.298 | 0.289  |            | -3.02  | 50    |
| Carbon Tetrachloride           | 0.494 | 0.480  | 0.1        | -2.83  | 50    |
| 2,2-Dichloropropane            | 0.662 | 0.630  |            | -4.83  | 50    |
| cis-1,2-Dichloroethene         | 2.866 | 2.731  |            | -4.71  | 50    |
| Chloroform                     | 4.355 | 4.229  | 0.2        | -2.89  | 50    |
| 1,1,1-Trichloroethane          | 0.618 | 0.613  | 0.1        | -0.81  | 50    |
| Methylcyclohexane              | 0.659 | 0.623  |            | -5.46  | 50    |
| 1,1-Dichloropropene            | 0.537 | 0.536  |            | -0.19  | 50    |
| Benzene                        | 1.710 | 1.702  | 0.5        | -0.47  | 50    |
| 1,2-Dichloroethane             | 0.535 | 0.527  | 0.1        | -1.5   | 50    |
| Trichloroethene                | 0.413 | 0.390  | 0.3        | -5.57  | 50    |
| 1,2-Dichloropropane            | 0.425 | 0.419  |            | -1.41  | 50    |
| Dibromomethane                 | 0.275 | 0.273  |            | -0.73  | 50    |
| Bromodichloromethane           | 0.590 | 0.577  | 0.2        | -2.2   | 50    |
| 4-Methyl-2-Pentanone           | 0.681 | 0.675  |            | -0.88  | 50    |
| Toluene                        | 2.012 | 2.024  | 0.4        | 0.6    | 50    |
| t-1,3-Dichloropropene          | 0.667 | 0.639  | 0.1        | -4.2   | 50    |
| cis-1,3-Dichloropropene        | 0.706 | 0.700  | 0.2        | -0.85  | 50    |
| 1,1,2-Trichloroethane          | 0.385 | 0.372  | 0.1        | -3.38  | 50    |
| 1,3-Dichloropropane            | 0.687 | 0.678  |            | -1.31  | 50    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.

VOLATILE CONTINUING CALIBRATION CHECK

Lab Name: CHEMTECH Contract: TETR06

Lab Code: CHEM Case No.: Q1774 SAS No.: Q1774 SDG No.: Q1774

Instrument ID: MSVOA\_N Calibration Date/Time: 04/11/2025 17:28

Lab File ID: VN086255.D Init. Calib. Date(s): 04/11/2025 04/11/2025

Heated Purge: (Y/N) N Init. Calib. Time(s): 11:04 12:40

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

| COMPOUND                    | RRF   | RRF020 | MIN<br>RRF | %D     | MAX%D |
|-----------------------------|-------|--------|------------|--------|-------|
| 2-Hexanone                  | 0.507 | 0.497  |            | -1.97  | 50    |
| Dibromochloromethane        | 0.422 | 0.409  | 0.1        | -3.08  | 50    |
| 1,2-Dibromoethane           | 0.389 | 0.381  |            | -2.06  | 50    |
| Tetrachloroethene           | 0.443 | 0.436  | 0.2        | -1.58  | 50    |
| Chlorobenzene               | 1.229 | 1.213  | 0.5        | -1.3   | 50    |
| 1,1,1,2-Tetrachloroethane   | 0.406 | 0.409  |            | 0.74   | 50    |
| Hexachloroethane            | 0.281 | 0.270  |            | -3.91  | 50    |
| Ethyl Benzene               | 2.232 | 2.191  | 0.1        | -1.84  | 50    |
| m/p-Xylenes                 | 0.830 | 0.800  | 0.3        | -3.61  | 50    |
| o-Xylene                    | 0.816 | 0.819  | 0.3        | 0.37   | 50    |
| Styrene                     | 1.376 | 1.320  | 0.3        | -4.07  | 50    |
| Bromoform                   | 0.276 | 0.256  | 0.1        | -7.25  | 50    |
| Isopropylbenzene            | 2.012 | 1.938  |            | -3.68  | 50    |
| 1,1,2,2-Tetrachloroethane   | 0.573 | 0.586  | 0.3        | 2.27   | 50    |
| 1,2,3-Trichloropropane      | 0.596 | 0.572  |            | -4.03  | 50    |
| Bromobenzene                | 0.446 | 0.429  |            | -3.81  | 50    |
| n-propylbenzene             | 2.376 | 2.251  |            | -5.26  | 50    |
| 2-Chlorotoluene             | 1.494 | 1.457  |            | -2.48  | 50    |
| 1,3,5-Trimethylbenzene      | 1.653 | 1.573  |            | -4.84  | 50    |
| 4-Chlorotoluene             | 1.492 | 1.416  |            | -5.09  | 50    |
| tert-butylbenzene           | 1.430 | 1.362  |            | -4.76  | 50    |
| 1,2,4-Trimethylbenzene      | 1.676 | 1.585  |            | -5.43  | 50    |
| sec-Butylbenzene            | 1.995 | 1.884  |            | -5.56  | 50    |
| p-Isopropyltoluene          | 1.664 | 1.531  |            | -7.99  | 50    |
| 1,3-Dichlorobenzene         | 0.829 | 0.766  | 0.2        | -7.6   | 50    |
| 1,4-Dichlorobenzene         | 0.831 | 0.752  | 0.2        | -9.51  | 50    |
| n-Butylbenzene              | 1.502 | 1.334  |            | -11.19 | 50    |
| 1,2-Dichlorobenzene         | 0.818 | 0.769  | 0.2        | -5.99  | 50    |
| 1,2-Dibromo-3-Chloropropane | 0.130 | 0.116  |            | -10.77 | 50    |
| 1,2,4-Trichlorobenzene      | 0.398 | 0.331  | 0.2        | -16.83 | 50    |
| Hexachlorobutadiene         | 0.172 | 0.126  |            | -26.74 | 50    |
| Naphthalene                 | 1.569 | 1.309  |            | -16.57 | 50    |
| 1,2,3-Trichlorobenzene      | 0.398 | 0.331  |            | -16.83 | 50    |
| 1,2-Dichloroethane-d4       | 2.926 | 2.837  | 0.01       | -3.04  | 50    |
| Toluene-d8                  | 1.694 | 1.736  | 0.01       | 2.48   | 50    |
| 4-Bromofluorobenzene        | 0.597 | 0.593  | 0.2        | -0.67  | 50    |
| Methyl Iodide               | 2.626 | 2.513  |            | -4.3   | 50    |
| Methyl methacrylate         | 0.491 | 0.468  |            | -4.68  | 50    |

All other compounds must meet a minimum RRF of 0.010.  
RRF of 1,4-Dioxane = Value should be divide by 1000.