

VOLATILE CONTINUING CALIBRATION CHECK

|                     |                   |                        |                              |
|---------------------|-------------------|------------------------|------------------------------|
| Lab Name:           | <u>Alliance</u>   | Contract:              | <u>NOBI03</u>                |
| Lab Code:           | <u>ACE</u>        | SDG No.:               | <u>Q2638</u>                 |
| Instrument ID:      | <u>MSVOA_Y</u>    | Calibration Date/Time: | <u>07/18/2025 21:30</u>      |
| Lab File ID:        | <u>VY023013.D</u> | Init. Calib. Date(s):  | <u>07/18/2025 07/18/2025</u> |
| Heated Purge: (Y/N) | <u>Y</u>          | Init. Calib. Time(s):  | <u>10:08 12:49</u>           |
| GC Column:          | <u>RXI-624</u>    | ID:                    | <u>0.25</u> (mm)             |

| COMPOUND                       | RRF   | RRF050 | MIN<br>RRF | %D     | MAX%D |
|--------------------------------|-------|--------|------------|--------|-------|
| Dichlorodifluoromethane        | 0.425 | 0.373  |            | -12.23 | 50    |
| Chloromethane                  | 0.722 | 0.855  | 0.1        | 18.42  | 50    |
| Vinyl Chloride                 | 0.915 | 1.043  |            | 13.99  | 50    |
| Bromomethane                   | 0.660 | 0.825  |            | 25     | 50    |
| Chloroethane                   | 0.611 | 0.741  |            | 21.28  | 50    |
| Tetrahydrofuran                | 0.083 | 0.111  |            | 33.74  | 50    |
| Trichlorofluoromethane         | 1.062 | 1.154  |            | 8.66   | 50    |
| 1,1,2-Trichlorotrifluoroethane | 0.558 | 0.528  |            | -5.38  | 50    |
| 1,1-Dichloroethene             | 0.531 | 0.499  |            | -6.03  | 50    |
| Acrylonitrile                  | 0.105 | 0.131  |            | 24.76  | 50    |
| Acetone                        | 0.092 | 0.096  |            | 4.35   | 50    |
| Carbon Disulfide               | 1.610 | 1.537  |            | -4.53  | 50    |
| Methyl tert-butyl Ether        | 1.274 | 1.467  |            | 15.15  | 50    |
| Methylene Chloride             | 0.871 | 0.702  |            | -19.4  | 50    |
| trans-1,2-Dichloroethene       | 0.591 | 0.581  |            | -1.69  | 50    |
| 1,1-Dichloroethane             | 1.079 | 1.114  | 0.1        | 3.24   | 50    |
| 2-Butanone                     | 0.132 | 0.164  |            | 24.24  | 50    |
| Carbon Tetrachloride           | 0.543 | 0.506  |            | -6.81  | 50    |
| 2,2-Dichloropropane            | 0.945 | 0.836  |            | -11.53 | 50    |
| cis-1,2-Dichloroethene         | 0.682 | 0.694  |            | 1.76   | 50    |
| Chloroform                     | 1.088 | 1.144  |            | 5.15   | 50    |
| 1,1,1-Trichloroethane          | 0.987 | 0.972  |            | -1.52  | 50    |
| 1,1-Dichloropropene            | 0.500 | 0.481  |            | -3.8   | 50    |
| Benzene                        | 1.504 | 1.533  |            | 1.93   | 50    |
| 1,2-Dichloroethane             | 0.374 | 0.421  |            | 12.57  | 50    |
| Trichloroethene                | 0.386 | 0.373  |            | -3.37  | 50    |
| 1,2-Dichloropropane            | 0.351 | 0.362  |            | 3.13   | 50    |
| Dibromomethane                 | 0.182 | 0.202  |            | 10.99  | 50    |
| Bromodichloromethane           | 0.502 | 0.528  |            | 5.18   | 50    |
| 4-Methyl-2-Pentanone           | 0.187 | 0.249  |            | 33.15  | 50    |
| Toluene                        | 0.935 | 0.953  |            | 1.92   | 50    |
| t-1,3-Dichloropropene          | 0.431 | 0.465  |            | 7.89   | 50    |
| cis-1,3-Dichloropropene        | 0.519 | 0.536  |            | 3.28   | 50    |
| 1,1,2-Trichloroethane          | 0.239 | 0.272  |            | 13.81  | 50    |
| 1,3-Dichloropropane            | 0.412 | 0.471  |            | 14.32  | 50    |
| 2-Hexanone                     | 0.123 | 0.163  |            | 32.52  | 50    |
| Dibromochloromethane           | 0.312 | 0.352  |            | 12.82  | 50    |
| 1,2-Dibromoethane              | 0.217 | 0.253  |            | 16.59  | 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: Alliance  
Lab Code: ACE  
Instrument ID: MSVOA\_Y  
Lab File ID: VY023013.D  
Heated Purge: (Y/N) Y  
GC Column: RXI-624 ID: 0.25 (mm)

Contract: NOBI03  
SDG No.: Q2638  
Calibration Date/Time: 07/18/2025 21:30  
Init. Calib. Date(s): 07/18/2025 07/18/2025  
Init. Calib. Time(s): 10:08 12:49

| COMPOUND                    | RRF   | RRF050 | MIN<br>RRF | %D     | MAX%D |
|-----------------------------|-------|--------|------------|--------|-------|
| Tetrachloroethene           | 0.511 | 0.488  |            | -4.5   | 50    |
| Chlorobenzene               | 1.180 | 1.137  | 0.3        | -3.64  | 50    |
| 1,1,1,2-Tetrachloroethane   | 0.402 | 0.396  |            | -1.49  | 50    |
| Ethyl Benzene               | 2.091 | 2.011  |            | -3.83  | 50    |
| m/p-Xylenes                 | 0.802 | 0.770  |            | -3.99  | 50    |
| o-Xylene                    | 0.744 | 0.725  |            | -2.55  | 50    |
| Styrene                     | 1.219 | 1.247  |            | 2.3    | 50    |
| Bromoform                   | 0.204 | 0.223  | 0.1        | 9.31   | 50    |
| Isopropylbenzene            | 4.239 | 3.692  |            | -12.9  | 50    |
| 1,1,2,2-Tetrachloroethane   | 0.613 | 0.642  | 0.3        | 4.73   | 50    |
| 1,2,3-Trichloropropane      | 0.623 | 0.602  |            | -3.37  | 50    |
| Bromobenzene                | 0.919 | 0.852  |            | -7.29  | 50    |
| n-propylbenzene             | 5.048 | 4.398  |            | -12.88 | 50    |
| 2-Chlorotoluene             | 2.839 | 2.531  |            | -10.85 | 50    |
| 1,3,5-Trimethylbenzene      | 3.343 | 2.992  |            | -10.5  | 50    |
| 4-Chlorotoluene             | 2.864 | 2.624  |            | -8.38  | 50    |
| tert-Butylbenzene           | 3.020 | 2.640  |            | -12.58 | 50    |
| 1,2,4-Trimethylbenzene      | 3.291 | 3.002  |            | -8.78  | 50    |
| sec-Butylbenzene            | 4.526 | 3.873  |            | -14.43 | 50    |
| p-Isopropyltoluene          | 3.676 | 3.177  |            | -13.57 | 50    |
| 1,3-Dichlorobenzene         | 1.855 | 1.691  |            | -8.84  | 50    |
| 1,4-Dichlorobenzene         | 1.795 | 1.636  |            | -8.86  | 50    |
| n-Butylbenzene              | 3.392 | 2.844  |            | -16.16 | 50    |
| 1,2-Dichlorobenzene         | 1.565 | 1.494  |            | -4.54  | 50    |
| 1,2-Dibromo-3-Chloropropane | 0.091 | 0.107  |            | 17.58  | 50    |
| 1,2,4-Trichlorobenzene      | 0.825 | 0.762  |            | -7.64  | 50    |
| Hexachlorobutadiene         | 0.520 | 0.384  |            | -26.15 | 50    |
| 1,2,3-Trichlorobenzene      | 0.681 | 0.669  |            | -1.76  | 50    |
| 1,2-Dichloroethane-d4       | 0.506 | 0.604  |            | 19.37  | 50    |
| Dibromofluoromethane        | 0.317 | 0.334  |            | 5.36   | 50    |
| Toluene-d8                  | 1.258 | 1.289  |            | 2.46   | 50    |
| 4-Bromofluorobenzene        | 0.396 | 0.418  |            | 5.56   | 50    |
| trans-1,4-Dichloro-2-butene | 0.209 | 0.209  |            | 0      | 50    |

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