

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU021325\  
 Data File : VU063262.D  
 Acq On : 13 Feb 2025 08:36  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Feb 14 04:40:50 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U021025DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Feb 11 08:42:19 2025  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

|       | Compound                    | Amount  | Calc.  | %Dev    | Area% | Dev(min) |
|-------|-----------------------------|---------|--------|---------|-------|----------|
| 1 i   | Fluorobenzene               | 1.000   | 1.000  | 0.0     | 93    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 9.970  | 0.3     | 99    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 9.446  | 5.5     | 94    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 9.939  | 0.6     | 94    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 11.198 | -12.0   | 99    | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 9.468  | 5.3     | 93    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 10.001 | -0.0    | 97    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 10.385 | -3.8    | 105   | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 10.108 | -1.1    | 97    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 10.372 | -3.7    | 95    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 10.258 | -2.6    | 96    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 21.419 | -7.1    | 90    | 0.00     |
| 13 T  | Acetone                     | 50.000  | 51.212 | -2.4    | 91    | 0.00     |
| 14 T  | Carbon Disulfide            | 10.000  | 9.851  | 1.5     | 95    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 10.078 | -0.8    | 97    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 10.269 | -2.7    | 97    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 10.133 | -1.3    | 96    | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 53.920 | -7.8    | 92    | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 11.000 | -10.0   | 99    | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 11.488 | -14.9   | 104   | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 10.502 | -5.0    | 100   | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 10.678 | -6.8    | 99    | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 9.774  | 2.3     | 92    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 97.059 | 2.9     | 86    | 0.00     |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 10.626 | -6.3    | 94    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 10.298 | -3.0    | 96    | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 10.226 | -2.3    | 96    | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 10.367 | -3.7    | 98    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 10.882 | -8.8    | 100   | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 10.498 | -5.0    | 99    | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 10.792 | -7.9    | 97    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 10.965 | -9.6    | 96    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 11.169 | -11.7   | 94    | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 55.852 | -11.7   | 91    | 0.00     |
| 35 RT | Benzene                     | 10.000  | 10.417 | -4.2    | 97    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 10.147 | -1.5    | 95    | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 10.460 | -4.6    | 98    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 10.331 | -3.3    | 95    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 11.811 | -18.1   | 91    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 10.826 | -8.3    | 91    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 20.669 | -3.3    | 88    | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 10.786 | -7.9    | 98    | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 10.154 | -1.5    | 95    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.646 | -6.5    | 97    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 55.115 | -10.2   | 89    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 47.833 | -139.2# | 213   | -0.01    |
| 47 t  | Methyl methacrylate         | 20.000  | 22.854 | -14.3   | 91    | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 11.777 | -17.8   | 90    | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 11.073 | -10.7   | 98    | 0.00     |

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 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Feb 14 04:40:50 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U021025DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Feb 11 08:42:19 2025  
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min  
 Max. RRF Dev : 30% Max. Rel. Area : 200%

|       | Compound                    | Amount | Calc.  | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|--------|--------|-------|-------|----------|
| 50 T  | t-1,3-Dichloropropene       | 10.000 | 11.433 | -14.3 | 97    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 10.974 | -9.7  | 96    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 10.409 | -4.1  | 94    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 10.448 | -4.5  | 94    | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 54.661 | -9.3  | 88    | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.851 | -8.5  | 96    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 10.494 | -4.9  | 92    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.288  | -28.8 | 111   | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.612 | -6.1  | 102   | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.822 | -8.2  | 98    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 10.720 | -7.2  | 98    | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 10.465 | -4.6  | 97    | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.970 | -9.7  | 98    | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 11.412 | -14.1 | 98    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 23.502 | -17.5 | 99    | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 11.342 | -13.4 | 97    | 0.00     |
| 66 RT | Styrene                     | 10.000 | 11.887 | -18.9 | 97    | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 10.863 | -8.6  | 93    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 1.057  | -5.7  | 99    | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 11.602 | -16.0 | 99    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 10.469 | -4.7  | 94    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 10.245 | -2.4  | 104   | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 10.881 | -8.8  | 96    | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 12.000 | -20.0 | 100   | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 11.545 | -15.4 | 100   | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 11.906 | -19.1 | 99    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 11.511 | -15.1 | 100   | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 11.330 | -13.3 | 98    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 12.019 | -20.2 | 99    | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 11.798 | -18.0 | 101   | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 47.324 | 5.4   | 87    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 12.067 | -20.7 | 101   | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 11.119 | -11.2 | 101   | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 11.131 | -11.3 | 98    | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 12.498 | -25.0 | 104   | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 10.882 | -8.8  | 98    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 10.704 | -7.0  | 85    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 12.188 | -21.9 | 100   | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 11.369 | -13.7 | 113   | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 9.007  | 9.9   | 85    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 11.635 | -16.3 | 95    | 0.00     |

(#) = Out of Range

SPCC's out = 0 CCC's out = 0