

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU110525\  
 Data File : VU063667.D  
 Acq On : 05 Nov 2025 10:13  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 5.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Nov 06 03:58:35 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U103125DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Mon Nov 03 00:51:51 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  | 98    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 10.170 | -1.7 | 96    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 9.800  | 2.0  | 97    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 10.247 | -2.5 | 95    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 9.813  | 1.9  | 100   | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 10.455 | -4.6 | 99    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 10.554 | -5.5 | 98    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 10.632 | -6.3 | 100   | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 10.567 | -5.7 | 99    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 10.494 | -4.9 | 95    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 10.178 | -1.8 | 95    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 18.539 | 7.3  | 82    | 0.00     |
| 13 T  | Acetone                     | 50.000  | 48.905 | 2.2  | 96    | 0.00     |
| 14 T  | Carbon Disulfide            | 10.000  | 10.186 | -1.9 | 97    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 9.271  | 7.3  | 97    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 10.461 | -4.6 | 97    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 10.421 | -4.2 | 98    | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 47.648 | 4.7  | 90    | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 10.618 | -6.2 | 97    | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 10.406 | -4.1 | 97    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 10.538 | -5.4 | 99    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 10.518 | -5.2 | 97    | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 10.060 | -0.6 | 91    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 72.705 | 27.3 | 74    | 0.00     |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 9.808  | 1.9  | 90    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 9.969  | 0.3  | 93    | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 10.521 | -5.2 | 100   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 10.473 | -4.7 | 98    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 10.645 | -6.4 | 98    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 10.910 | -9.1 | 99    | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 10.135 | -1.3 | 94    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 9.829  | 1.7  | 93    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 9.998  | 0.0  | 93    | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 45.605 | 8.8  | 80    | 0.00     |
| 35 RT | Benzene                     | 10.000  | 10.441 | -4.4 | 98    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 10.144 | -1.4 | 93    | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 10.835 | -8.4 | 107   | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 10.277 | -2.8 | 96    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 8.704  | 13.0 | 87    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 9.179  | 8.2  | 91    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 18.879 | 5.6  | 82    | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 10.876 | -8.8 | 97    | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 9.970  | 0.3  | 90    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.112 | -1.1 | 92    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 46.023 | 8.0  | 82    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 17.733 | 11.3 | 78    | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 18.993 | 5.0  | 83    | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 9.693  | 3.1  | 84    | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 10.423 | -4.2 | 96    | 0.00     |

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 Sample : VSTDCCC010  
 Misc : 5.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Nov 06 03:58:35 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U103125DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Mon Nov 03 00:51:51 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 | 9.483  | 5.2   | 88    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 9.879  | 1.2   | 91    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 10.048 | -0.5  | 91    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 9.765  | 2.3   | 91    | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 46.848 | 6.3   | 85    | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.068 | -0.7  | 90    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 9.517  | 4.8   | 88    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.148  | -14.8 | 107   | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.927 | -9.3  | 99    | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.486 | -4.9  | 98    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 10.194 | -1.9  | 96    | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 9.630  | 3.7   | 91    | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.196 | -2.0  | 91    | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 10.298 | -3.0  | 96    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 20.920 | -4.6  | 97    | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 10.372 | -3.7  | 95    | 0.00     |
| 66 RT | Styrene                     | 10.000 | 10.307 | -3.1  | 94    | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 9.799  | 2.0   | 88    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 0.956  | 4.4   | 84    | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 10.433 | -4.3  | 96    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 9.277  | 7.2   | 84    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 9.241  | 7.6   | 91    | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 10.345 | -3.5  | 95    | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 10.445 | -4.5  | 95    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 10.417 | -4.2  | 97    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 10.272 | -2.7  | 95    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 10.317 | -3.2  | 95    | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 10.102 | -1.0  | 93    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 10.464 | -4.6  | 93    | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 10.289 | -2.9  | 94    | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 26.078 | 47.8# | 50    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 10.265 | -2.7  | 93    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 10.357 | -3.6  | 94    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 10.386 | -3.9  | 95    | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 10.238 | -2.4  | 90    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 10.242 | -2.4  | 94    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 9.221  | 7.8   | 78    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 10.020 | -0.2  | 87    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 10.857 | -8.6  | 98    | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 8.362  | 16.4  | 74    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 9.762  | 2.4   | 84    | 0.00     |

(#) = Out of Range

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