

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU072823\  
 Data File : VU054905.D  
 Acq On : 28 Jul 2023 10:05  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Jul 28 23:06:05 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U072523DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Wed Jul 26 09:45:35 2023  
 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  | 9.631  | 3.7   | 97    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 9.379  | 6.2   | 96    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 9.438  | 5.6   | 95    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 8.922  | 10.8  | 89    | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 8.946  | 10.5  | 85    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 8.207  | 17.9  | 85    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 9.509  | 4.9   | 99    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 9.234  | 7.7   | 92    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 9.666  | 3.3   | 93    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 9.598  | 4.0   | 98    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 19.171 | 4.1   | 95    | 0.00     |
| 13 T  | Acetone                     | 50.000  | 56.612 | -13.2 | 116   | 0.00     |
| 14 T  | Carbon Disulfide            | 10.000  | 8.651  | 13.5  | 92    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 10.087 | -0.9  | 104   | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 9.847  | 1.5   | 100   | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 9.789  | 2.1   | 100   | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 56.125 | -12.3 | 104   | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 10.416 | -4.2  | 104   | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 9.532  | 4.7   | 94    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 9.464  | 5.4   | 100   | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 10.175 | -1.8  | 101   | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 8.944  | 10.6  | 91    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 95.867 | 4.1   | 116   | -0.03    |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 9.737  | 2.6   | 96    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 9.602  | 4.0   | 96    | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 10.067 | -0.7  | 100   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 9.589  | 4.1   | 97    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 9.355  | 6.4   | 94    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 9.869  | 1.3   | 101   | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 9.821  | 1.8   | 98    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 9.613  | 3.9   | 97    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 9.611  | 3.9   | 97    | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 55.532 | -11.1 | 96    | 0.00     |
| 35 RT | Benzene                     | 10.000  | 9.722  | 2.8   | 97    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 9.940  | 0.6   | 100   | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 9.749  | 2.5   | 97    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 10.070 | -0.7  | 98    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 10.219 | -2.2  | 97    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 11.008 | -10.1 | 98    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 18.362 | 8.2   | 98    | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 9.531  | 4.7   | 100   | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 10.172 | -1.7  | 97    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.274 | -2.7  | 99    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 50.633 | -1.3  | 98    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 22.787 | -13.9 | 95    | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 20.927 | -4.6  | 97    | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 10.427 | -4.3  | 96    | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 9.968  | 0.3   | 99    | 0.00     |

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 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Jul 28 23:06:05 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U072523DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Wed Jul 26 09:45:35 2023  
 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 | 10.253 | -2.5  | 99    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 9.644  | 3.6   | 93    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 9.851  | 1.5   | 100   | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 10.188 | -1.9  | 101   | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 53.127 | -6.3  | 98    | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.039 | -0.4  | 97    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 10.305 | -3.0  | 98    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.279  | -27.9 | 119   | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.194 | -1.9  | 103   | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 9.783  | 2.2   | 98    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 9.989  | 0.1   | 98    | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 9.910  | 0.9   | 93    | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.638 | -6.4  | 98    | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 10.093 | -0.9  | 98    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 20.530 | -2.7  | 100   | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 10.057 | -0.6  | 98    | 0.00     |
| 66 RT | Styrene                     | 10.000 | 10.274 | -2.7  | 97    | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 11.201 | -12.0 | 98    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 1.065  | -6.5  | 107   | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 10.158 | -1.6  | 99    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 10.273 | -2.7  | 97    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 9.689  | 3.1   | 98    | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 10.308 | -3.1  | 98    | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 10.341 | -3.4  | 98    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 10.329 | -3.3  | 99    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 10.430 | -4.3  | 99    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 10.364 | -3.6  | 96    | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 10.221 | -2.2  | 98    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 10.419 | -4.2  | 98    | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 10.501 | -5.0  | 99    | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 54.177 | -8.4  | 95    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 10.726 | -7.3  | 98    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 10.071 | -0.7  | 99    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 10.257 | -2.6  | 100   | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 10.527 | -5.3  | 91    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 9.884  | 1.2   | 99    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 10.372 | -3.7  | 92    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 9.848  | 1.5   | 94    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 10.295 | -2.9  | 98    | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 9.065  | 9.4   | 86    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 9.598  | 4.0   | 94    | 0.00     |

(#) = Out of Range

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