

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU080624\  
 Data File : VU060314.D  
 Acq On : 06 Aug 2024 09:55  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Aug 07 06:46:32 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U080524DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Aug 06 11:07:28 2024  
 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  | 105   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 9.712  | 2.9  | 98    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 9.053  | 9.5  | 96    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 9.337  | 6.6  | 95    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 8.830  | 11.7 | 91    | -0.02    |
| 6 T   | Chloroethane                | 10.000  | 9.399  | 6.0  | 96    | -0.01    |
| 7 T   | Trichlorofluoromethane      | 10.000  | 9.698  | 3.0  | 98    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 9.636  | 3.6  | 98    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 9.543  | 4.6  | 99    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 9.499  | 5.0  | 97    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 9.557  | 4.4  | 98    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 17.583 | 12.1 | 97    | -0.01    |
| 13 T  | Acetone                     | 50.000  | 50.940 | -1.9 | 141   | -0.02    |
| 14 T  | Carbon Disulfide            | 10.000  | 9.367  | 6.3  | 96    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 9.146  | 8.5  | 98    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 9.507  | 4.9  | 97    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 9.449  | 5.5  | 98    | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 49.145 | 1.7  | 118   | -0.02    |
| 19    | Cyclohexane                 | 10.000  | 9.695  | 3.0  | 97    | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 10.778 | -7.8 | 96    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 9.805  | 2.0  | 103   | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 9.725  | 2.8  | 99    | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 9.171  | 8.3  | 97    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 82.837 | 17.2 | 97    | -0.05    |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 9.204  | 8.0  | 97    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 9.642  | 3.6  | 100   | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 9.505  | 4.9  | 100   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 9.642  | 3.6  | 100   | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 9.661  | 3.4  | 99    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 9.955  | 0.4  | 101   | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 9.431  | 5.7  | 97    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 9.393  | 6.1  | 97    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 9.966  | 0.3  | 98    | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 41.196 | 17.6 | 96    | -0.02    |
| 35 RT | Benzene                     | 10.000  | 9.891  | 1.1  | 98    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 9.450  | 5.5  | 99    | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 9.744  | 2.6  | 97    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 10.030 | -0.3 | 99    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 8.408  | 15.9 | 96    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 8.308  | 16.9 | 95    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 15.401 | 23.0 | 94    | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 9.799  | 2.0  | 99    | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 9.307  | 6.9  | 96    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 9.920  | 0.8  | 101   | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 49.896 | 0.2  | 99    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 21.557 | -7.8 | 107   | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 21.040 | -5.2 | 98    | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 10.741 | -7.4 | 98    | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 10.923 | -9.2 | 98    | 0.00     |

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 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Aug 07 06:46:32 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U080524DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Aug 06 11:07:28 2024  
 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 | 10.422 | -4.2   | 99    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 10.079 | -0.8   | 97    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 9.549  | 4.5    | 99    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 9.909  | 0.9    | 99    | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 59.305 | -18.6  | 119   | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.073 | -0.7   | 101   | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 9.592  | 4.1    | 97    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.366  | -36.6# | 128   | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.461 | -4.6   | 103   | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.685 | -6.9   | 101   | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 10.042 | -0.4   | 101   | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 10.407 | -4.1   | 103   | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.672 | -6.7   | 104   | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 9.549  | 4.5    | 99    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 19.836 | 0.8    | 101   | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 9.646  | 3.5    | 100   | 0.00     |
| 66 RT | Styrene                     | 10.000 | 9.978  | 0.2    | 102   | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 10.120 | -1.2   | 105   | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 1.056  | -5.6   | 104   | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 9.764  | 2.4    | 101   | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 9.748  | 2.5    | 101   | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 9.587  | 4.1    | 91    | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 11.105 | -11.1  | 102   | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 10.013 | -0.1   | 102   | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 10.061 | -0.6   | 101   | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 10.127 | -1.3   | 103   | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 10.163 | -1.6   | 102   | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 9.824  | 1.8    | 102   | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 10.144 | -1.4   | 104   | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 10.063 | -0.6   | 103   | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 44.863 | 10.3   | 93    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 10.016 | -0.2   | 102   | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 11.329 | -13.3  | 103   | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 11.692 | -16.9  | 105   | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 10.007 | -0.1   | 102   | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 11.305 | -13.0  | 105   | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 9.659  | 3.4    | 100   | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 10.692 | -6.9   | 96    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 11.006 | -10.1  | 102   | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 9.927  | 0.7    | 92    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 10.778 | -7.8   | 96    | 0.00     |

(#) = Out of Range

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