

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU031721\  
 Data File : VU042698.D  
 Acq On : 17 Mar 2021 19:37  
 Operator : SY/MD  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Mar 18 07:38:11 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U031721DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Mar 18 07:25:09 2021  
 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   | 101   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 10.327 | -3.3  | 105   | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 9.688  | 3.1   | 104   | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 10.075 | -0.7  | 104   | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 10.335 | -3.4  | 101   | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 11.098 | -11.0 | 112   | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 10.929 | -9.3  | 110   | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 11.030 | -10.3 | 112   | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 10.556 | -5.6  | 109   | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 11.322 | -13.2 | 112   | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 10.513 | -5.1  | 110   | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 21.455 | -7.3  | 107   | 0.00     |
| 13 T  | Acetone                     | 50.000  | 46.890 | 6.2   | 95    | 0.00     |
| 14 T  | Carbon Disulfide            | 10.000  | 10.387 | -3.9  | 107   | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 9.873  | 1.3   | 107   | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 10.438 | -4.4  | 109   | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 10.187 | -1.9  | 106   | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 46.795 | 6.4   | 96    | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 10.876 | -8.8  | 112   | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 11.189 | -11.9 | 107   | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 9.675  | 3.2   | 101   | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 10.166 | -1.7  | 107   | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 10.822 | -8.2  | 109   | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 74.554 | 25.4  | 63    | 0.00     |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 10.463 | -4.6  | 110   | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 10.297 | -3.0  | 105   | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 10.132 | -1.3  | 106   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 10.134 | -1.3  | 104   | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 10.463 | -4.6  | 105   | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 10.596 | -6.0  | 105   | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 10.325 | -3.2  | 108   | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 10.158 | -1.6  | 105   | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 10.722 | -7.2  | 105   | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 54.939 | -9.9  | 96    | 0.00     |
| 35 RT | Benzene                     | 10.000  | 10.352 | -3.5  | 105   | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 10.247 | -2.5  | 107   | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 10.767 | -7.7  | 108   | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 10.461 | -4.6  | 107   | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 9.541  | 4.6   | 99    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 10.382 | -3.8  | 97    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 18.429 | 7.9   | 98    | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 10.392 | -3.9  | 108   | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 10.706 | -7.1  | 103   | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.623 | -6.2  | 105   | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 55.357 | -10.7 | 105   | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 16.365 | 18.2  | 87    | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 22.140 | -10.7 | 104   | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 11.426 | -14.3 | 106   | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 11.303 | -13.0 | 106   | 0.00     |

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 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 16 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Mar 18 07:38:11 2021  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U031721DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Mar 18 07:25:09 2021  
 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 | 11.187 | -11.9 | 103   | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 10.734 | -7.3  | 103   | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 10.530 | -5.3  | 106   | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 10.622 | -6.2  | 105   | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 56.427 | -12.9 | 104   | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.921 | -9.2  | 105   | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 10.567 | -5.7  | 106   | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.060  | -6.0  | 97    | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.519 | -5.2  | 107   | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.933 | -9.3  | 106   | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 11.006 | -10.1 | 105   | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 11.451 | -14.5 | 108   | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 11.809 | -18.1 | 107   | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 11.840 | -18.4 | 107   | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 23.969 | -19.8 | 108   | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 11.844 | -18.4 | 108   | 0.00     |
| 66 RT | Styrene                     | 10.000 | 12.380 | -23.8 | 109   | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 11.254 | -12.5 | 103   | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 1.120  | -12.0 | 105   | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 12.059 | -20.6 | 109   | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 11.370 | -13.7 | 108   | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 12.498 | -25.0 | 118   | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 11.310 | -13.1 | 108   | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 11.965 | -19.6 | 106   | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 11.566 | -15.7 | 106   | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 12.279 | -22.8 | 107   | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 12.098 | -21.0 | 108   | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 11.832 | -18.3 | 106   | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 12.404 | -24.0 | 108   | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 11.981 | -19.8 | 106   | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 42.006 | 16.0  | 88    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 12.239 | -22.4 | 107   | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 11.246 | -12.5 | 105   | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 11.470 | -14.7 | 106   | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 12.382 | -23.8 | 107   | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 11.460 | -14.6 | 107   | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 11.602 | -16.0 | 101   | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 11.510 | -15.1 | 103   | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 11.117 | -11.2 | 104   | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 9.881  | 1.2   | 105   | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 11.630 | -16.3 | 103   | 0.00     |

(#) = Out of Range

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