

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU091622\  
 Data File : VU050727.D  
 Acq On : 16 Sep 2022 16:50  
 Operator : JC/MD  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Sep 17 01:08:24 2022  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U081922DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Aug 25 14:01:59 2022  
 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    | 91    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 10.219  | -2.2   | 90    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 7.865   | 21.3   | 71    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 9.006   | 9.9    | 78    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 10.965  | -9.6   | 102   | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 8.471   | 15.3   | 74    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 8.743   | 12.6   | 77    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 8.999   | 10.0   | 80    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 8.232   | 17.7   | 73    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 11.088  | -10.9  | 95    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 8.944   | 10.6   | 79    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 23.956  | -19.8  | 103   | 0.00     |
| 13 T  | Acetone                     | 50.000  | 59.977  | -20.0  | 107   | 0.00     |
| 14 T  | Carbon Disulfide            | 10.000  | 7.370   | 26.3   | 66    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 9.231   | 7.7    | 92    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 8.491   | 15.1   | 75    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 9.307   | 6.9    | 82    | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 67.267  | -34.5# | 114   | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 9.318   | 6.8    | 84    | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 11.395  | -13.9  | 87    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 8.907   | 10.9   | 84    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 9.463   | 5.4    | 84    | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 9.775   | 2.2    | 86    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 119.066 | -19.1  | 115   | 0.00     |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 10.217  | -2.2   | 91    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 11.733  | -17.3  | 107   | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 9.687   | 3.1    | 86    | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 9.267   | 7.3    | 82    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 9.621   | 3.8    | 86    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 9.614   | 3.9    | 83    | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 8.071   | 19.3   | 72    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 9.488   | 5.1    | 82    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 10.000  | 14.102  | -41.0# | 105   | 0.00     |
| 34 t  | Propionitrile               | 50.000  | 73.595  | -47.2# | 146   | 0.00     |
| 35 RT | Benzene                     | 10.000  | 10.407  | -4.1   | 89    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 11.001  | -10.0  | 97    | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 9.520   | 4.8    | 84    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 11.141  | -11.4  | 93    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 11.302  | -13.0  | 110   | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 12.248  | -22.5  | 120   | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 26.199  | -31.0# | 117   | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 9.758   | 2.4    | 89    | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 11.382  | -13.8  | 97    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.666  | -6.7   | 92    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 79.046  | -58.1# | 120   | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 18.897  | 5.5    | 80    | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 24.651  | -23.3  | 113   | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 11.173  | -11.7  | 103   | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 11.212  | -12.1  | 85    | 0.00     |

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 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

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 QLast Update : Thu Aug 25 14:01:59 2022  
 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 | 12.308 | -23.1  | 98    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 11.405 | -14.0  | 92    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 11.487 | -14.9  | 98    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 11.986 | -19.9  | 99    | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 79.562 | -59.1# | 122   | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.791 | -7.9   | 93    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 11.768 | -17.7  | 98    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.864  | -86.4# | 145   | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 9.126  | 8.7    | 79    | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.627 | -6.3   | 84    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 10.019 | -0.2   | 85    | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 10.429 | -4.3   | 88    | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.619 | -6.2   | 86    | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 8.861  | 11.4   | 80    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 18.524 | 7.4    | 82    | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 9.059  | 9.4    | 81    | 0.00     |
| 66 RT | Styrene                     | 10.000 | 9.097  | 9.0    | 81    | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 11.803 | -18.0  | 99    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 0.961  | 3.9    | 83    | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 8.889  | 11.1   | 80    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 12.508 | -25.1  | 107   | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 11.942 | -19.4  | 99    | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 10.373 | -3.7   | 82    | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 8.671  | 13.3   | 77    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 9.184  | 8.2    | 81    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 9.023  | 9.8    | 79    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 9.166  | 8.3    | 78    | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 8.802  | 12.0   | 79    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 8.772  | 12.3   | 78    | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 8.433  | 15.7   | 75    | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 54.544 | -9.1   | 94    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 7.941  | 20.6   | 70    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 10.257 | -2.6   | 80    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 10.313 | -3.1   | 80    | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 7.714  | 22.9   | 69    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 10.364 | -3.6   | 82    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 12.050 | -20.5  | 103   | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 9.498  | 5.0    | 72    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 8.410  | 15.9   | 69    | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 7.778  | 22.2   | 73    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 8.160  | 18.4   | 74    | 0.00     |

(#) = Out of Range

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