

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU073024\  
 Data File : VU060272.D  
 Acq On : 30 Jul 2024 19:27  
 Operator : MD/SY  
 Sample : VSTDICV010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 ICSVU073024

Quant Time: Jul 31 04:54:54 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U073024DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Wed Jul 31 04:49:33 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     | 92    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 10.000  | 5.177  | 48.2#   | 46    | 0.00     |
| 3 t   | Chloromethane               | 10.000  | 7.392  | 26.1    | 68    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 10.000  | 7.797  | 22.0    | 68    | 0.00     |
| 5 T   | Bromomethane                | 10.000  | 8.247  | 17.5    | 74    | 0.00     |
| 6 T   | Chloroethane                | 10.000  | 8.460  | 15.4    | 75    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 10.000  | 8.502  | 15.0    | 76    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 10.000  | 9.577  | 4.2     | 85    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 10.000  | 9.733  | 2.7     | 84    | 0.00     |
| 10 t  | Iodomethane                 | 10.000  | 9.609  | 3.9     | 83    | 0.00     |
| 11 t  | Allyl Chloride              | 10.000  | 10.326 | -3.3    | 93    | 0.00     |
| 12 t  | Acrylonitrile               | 20.000  | 49.197 | -146.0# | 212   | 0.00     |
| 13 T  | Acetone                     | 50.000  | 32.342 | 35.3#   | 63    | 0.01     |
| 14 T  | Carbon Disulfide            | 10.000  | 8.833  | 11.7    | 78    | 0.00     |
| 15 RT | Methylene Chloride          | 10.000  | 8.619  | 13.8    | 85    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 10.000  | 9.353  | 6.5     | 83    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 10.000  | 9.912  | 0.9     | 87    | 0.00     |
| 18 T  | 2-Butanone                  | 50.000  | 39.781 | 20.4    | 78    | 0.00     |
| 19    | Cyclohexane                 | 10.000  | 9.083  | 9.2     | 82    | 0.00     |
| 20    | Methylcyclohexane           | 10.000  | 9.857  | 1.4     | 78    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 10.000  | 8.832  | 11.7    | 79    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 10.000  | 10.264 | -2.6    | 90    | 0.00     |
| 23 t  | Diethyl Ether               | 10.000  | 9.134  | 8.7     | 80    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 100.000 | 50.086 | 49.9#   | 44    | 0.03     |
| 25 t  | Methyl tert-Butyl Ether     | 10.000  | 10.142 | -1.4    | 88    | 0.00     |
| 26 t  | Bromochloromethane          | 10.000  | 10.443 | -4.4    | 95    | 0.00     |
| 27 t  | Chloroform                  | 10.000  | 9.871  | 1.3     | 88    | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 10.000  | 9.671  | 3.3     | 85    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 10.000  | 9.358  | 6.4     | 81    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 10.000  | 9.595  | 4.0     | 84    | 0.00     |
| 31 t  | Isopropyl Ether             | 10.000  | 10.521 | -5.2    | 91    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 10.000  | 0.000  | 100.0#  | 0     | -4.50#   |
| 33    | Tert-Amyl methyl ether      | 10.000  | 0.000  | 100.0#  | 0     | -5.95#   |
| 34 t  | Propionitrile               | 50.000  | 98.219 | -96.4#  | 182   | 0.00     |
| 35 RT | Benzene                     | 10.000  | 10.034 | -0.3    | 85    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 10.000  | 9.286  | 7.1     | 82    | 0.00     |
| 37 RT | Trichloroethene             | 10.000  | 9.076  | 9.2     | 79    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 10.000  | 9.452  | 5.5     | 80    | 0.00     |
| 39 t  | Methacrylonitrile           | 10.000  | 9.627  | 3.7     | 88    | 0.00     |
| 40 t  | Methyl acrylate             | 10.000  | 10.001 | -0.0    | 89    | 0.00     |
| 41 t  | Tetrahydrofuran             | 20.000  | 48.132 | -140.7# | 218   | 0.00     |
| 42 t  | 1-Chlorobutane              | 10.000  | 9.877  | 1.2     | 86    | 0.00     |
| 43 T  | Dibromomethane              | 10.000  | 10.364 | -3.6    | 89    | 0.00     |
| 44 T  | Bromodichloromethane        | 10.000  | 10.290 | -2.9    | 89    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 50.000  | 56.184 | -12.4   | 88    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 20.000  | 10.762 | 46.2#   | 44    | 0.00     |
| 47 t  | Methyl methacrylate         | 20.000  | 9.819  | 50.9#   | 37    | 0.00     |
| 48 t  | Ethyl methacrylate          | 10.000  | 9.895  | 1.1     | 88    | 0.00     |
| 49 Rt | Toluene                     | 10.000  | 11.050 | -10.5   | 86    | 0.00     |

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 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 10 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 ICVVU073024

Quant Time: Jul 31 04:54:54 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U073024DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Wed Jul 31 04:49:33 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 | 10.255 | -2.6  | 83    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 10.000 | 9.842  | 1.6   | 81    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 10.000 | 9.785  | 2.1   | 84    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 10.000 | 10.045 | -0.4  | 83    | 0.00     |
| 54 t  | 2-Hexanone                  | 50.000 | 53.187 | -6.4  | 86    | 0.00     |
| 55 t  | Dibromochloromethane        | 10.000 | 10.489 | -4.9  | 89    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 10.000 | 10.228 | -2.3  | 86    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 1.000  | 1.145  | -14.5 | 95    | 0.00     |
| 58 RT | Tetrachloroethene           | 10.000 | 10.304 | -3.0  | 87    | 0.00     |
| 59 Rt | Chlorobenzene               | 10.000 | 10.775 | -7.8  | 87    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 10.000 | 9.705  | 2.9   | 83    | 0.00     |
| 61 t  | Pentachloroethane           | 10.000 | 10.317 | -3.2  | 89    | 0.00     |
| 62 t  | Hexachloroethane            | 10.000 | 10.576 | -5.8  | 89    | 0.00     |
| 63 Rt | Ethyl Benzene               | 10.000 | 9.666  | 3.3   | 85    | 0.00     |
| 64 RT | m/p-Xylenes                 | 20.000 | 19.788 | 1.1   | 88    | 0.00     |
| 65 RT | o-Xylene                    | 10.000 | 9.596  | 4.0   | 85    | 0.00     |
| 66 RT | Styrene                     | 10.000 | 10.205 | -2.1  | 91    | 0.00     |
| 67 t  | Bromoform                   | 10.000 | 10.461 | -4.6  | 89    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 1.000  | 1.113  | -11.3 | 101   | 0.00     |
| 69 T  | Isopropylbenzene            | 10.000 | 10.016 | -0.2  | 90    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 10.000 | 9.910  | 0.9   | 83    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 10.000 | 9.895  | 1.1   | 83    | 0.00     |
| 72 t  | Bromobenzene                | 10.000 | 11.159 | -11.6 | 91    | 0.00     |
| 73 t  | n-propylbenzene             | 10.000 | 9.754  | 2.5   | 87    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 10.000 | 11.622 | -16.2 | 89    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 10.000 | 10.000 | 0.0   | 90    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 10.000 | 10.101 | -1.0  | 90    | 0.00     |
| 77 t  | tert-Butylbenzene           | 10.000 | 9.986  | 0.1   | 90    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 10.000 | 9.895  | 1.1   | 88    | 0.00     |
| 79 t  | sec-Butylbenzene            | 10.000 | 9.881  | 1.2   | 88    | 0.00     |
| 80    | Nitrobenzene                | 50.000 | 11.294 | 77.4# | 19    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 10.000 | 9.994  | 0.1   | 89    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 10.000 | 11.122 | -11.2 | 90    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 10.000 | 11.309 | -13.1 | 89    | 0.00     |
| 84 t  | n-Butylbenzene              | 10.000 | 9.975  | 0.3   | 89    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 10.000 | 11.124 | -11.2 | 89    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 10.000 | 10.131 | -1.3  | 83    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 10.000 | 12.566 | -25.7 | 99    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 10.000 | 10.340 | -3.4  | 88    | 0.00     |
| 89 t  | Naphthalene                 | 10.000 | 10.422 | -4.2  | 96    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 10.000 | 10.177 | -1.8  | 91    | 0.00     |

(#) = Out of Range

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