

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU020224\  
 Data File : VU057784.D  
 Acq On : 02 Feb 2024 09:20  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Feb 03 05:50:53 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U013124DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Fri Feb 02 00:27:55 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|-------|-------|----------|
| 1 i   | Fluorobenzene               | 1.000 | 1.000 | 0.0   | 102   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.333 | 0.328 | 1.5   | 97    | 0.00     |
| 3 t   | Chloromethane               | 0.320 | 0.310 | 3.1   | 100   | 0.00     |
| 4 Rt  | Vinyl Chloride              | 0.326 | 0.331 | -1.5  | 99    | 0.00     |
| 5 T   | Bromomethane                | 0.193 | 0.194 | -0.5  | 108   | 0.00     |
| 6 T   | Chloroethane                | 0.206 | 0.211 | -2.4  | 103   | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.499 | 0.495 | 0.8   | 95    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 0.256 | 0.257 | -0.4  | 99    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 0.249 | 0.247 | 0.8   | 97    | 0.00     |
| 10 t  | Iodomethane                 | 0.295 | 0.320 | -8.5  | 98    | 0.00     |
| 11 t  | Allyl Chloride              | 0.337 | 0.355 | -5.3  | 100   | 0.00     |
| 12 t  | Acrylonitrile               | 0.042 | 0.048 | -14.3 | 102   | 0.00     |
| 13 T  | Acetone                     | 0.077 | 0.091 | -18.2 | 107   | 0.02     |
| 14 T  | Carbon Disulfide            | 0.795 | 0.809 | -1.8  | 99    | 0.00     |
| 15 RT | Methylene Chloride          | 0.297 | 0.279 | 6.1   | 98    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 0.271 | 0.279 | -3.0  | 98    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 0.512 | 0.535 | -4.5  | 102   | 0.00     |
| 18 T  | 2-Butanone                  | 0.083 | 0.098 | -18.1 | 107   | 0.01     |
| 19    | Cyclohexane                 | 0.401 | 0.402 | -0.2  | 98    | 0.00     |
| 20    | Methylcyclohexane           | 0.485 | 0.502 | -3.5  | 98    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 0.455 | 0.478 | -5.1  | 99    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 0.297 | 0.301 | -1.3  | 98    | 0.00     |
| 23 t  | Diethyl Ether               | 0.179 | 0.184 | -2.8  | 99    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 0.020 | 0.021 | -5.0  | 103   | 0.13     |
| 25 t  | Methyl tert-Butyl Ether     | 0.583 | 0.599 | -2.7  | 98    | 0.00     |
| 26 t  | Bromochloromethane          | 0.119 | 0.119 | 0.0   | 97    | 0.00     |
| 27 t  | Chloroform                  | 0.533 | 0.548 | -2.8  | 100   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 0.481 | 0.492 | -2.3  | 98    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 0.389 | 0.406 | -4.4  | 100   | 0.00     |
| 30 RT | Carbon Tetrachloride        | 0.414 | 0.426 | -2.9  | 98    | 0.00     |
| 31 t  | Isopropyl Ether             | 0.750 | 0.789 | -5.2  | 101   | 0.00     |
| 32    | Ethyl-t-butyl ether         | 0.696 | 0.726 | -4.3  | 96    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 0.586 | 0.597 | -1.9  | 94    | 0.00     |
| 34 t  | Propionitrile               | 0.017 | 0.019 | -11.8 | 98    | 0.02     |
| 35 RT | Benzene                     | 1.120 | 1.146 | -2.3  | 99    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 0.326 | 0.335 | -2.8  | 100   | 0.00     |
| 37 RT | Trichloroethene             | 0.297 | 0.301 | -1.3  | 98    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 0.290 | 0.299 | -3.1  | 100   | 0.00     |
| 39 t  | Methacrylonitrile           | 0.068 | 0.076 | -11.8 | 101   | 0.00     |
| 40 t  | Methyl acrylate             | 0.103 | 0.112 | -8.7  | 83    | 0.00     |
| 41 t  | Tetrahydrofuran             | 0.036 | 0.036 | 0.0   | 92    | 0.00     |
| 42 t  | 1-Chlorobutane              | 0.513 | 0.544 | -6.0  | 100   | 0.00     |
| 43 T  | Dibromomethane              | 0.134 | 0.139 | -3.7  | 99    | 0.00     |
| 44 T  | Bromodichloromethane        | 0.374 | 0.388 | -3.7  | 99    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 0.142 | 0.151 | -6.3  | 95    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 0.032 | 0.039 | -21.9 | 95    | 0.00     |
| 47 t  | Methyl methacrylate         | 0.118 | 0.126 | -6.8  | 98    | 0.00     |
| 48 t  | Ethyl methacrylate          | 0.236 | 0.244 | -3.4  | 92    | 0.00     |
| 49 Rt | Toluene                     | 0.684 | 0.711 | -3.9  | 99    | 0.00     |

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

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Feb 03 05:50:53 2024  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U013124DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Fri Feb 02 00:27:55 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                    | AvgRF | CCRF  | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|-------|-------|----------|
| 50 T  | t-1,3-Dichloropropene       | 0.323 | 0.348 | -7.7  | 98    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 0.409 | 0.433 | -5.9  | 99    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 0.184 | 0.188 | -2.2  | 97    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 0.331 | 0.338 | -2.1  | 100   | 0.00     |
| 54 t  | 2-Hexanone                  | 0.131 | 0.148 | -13.0 | 102   | 0.00     |
| 55 t  | Dibromochloromethane        | 0.228 | 0.234 | -2.6  | 97    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 0.169 | 0.177 | -4.7  | 97    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 0.379 | 0.479 | -26.4 | 119   | 0.00     |
| 58 RT | Tetrachloroethene           | 0.262 | 0.258 | 1.5   | 97    | 0.00     |
| 59 Rt | Chlorobenzene               | 0.757 | 0.770 | -1.7  | 95    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 0.257 | 0.265 | -3.1  | 96    | 0.00     |
| 61 t  | Pentachloroethane           | 0.187 | 0.206 | -10.2 | 94    | 0.00     |
| 62 t  | Hexachloroethane            | 0.204 | 0.220 | -7.8  | 90    | 0.00     |
| 63 Rt | Ethyl Benzene               | 1.345 | 1.401 | -4.2  | 96    | 0.00     |
| 64 RT | m/p-Xylenes                 | 0.489 | 0.516 | -5.5  | 95    | 0.00     |
| 65 RT | o-Xylene                    | 0.484 | 0.504 | -4.1  | 94    | 0.00     |
| 66 RT | Styrene                     | 0.722 | 0.782 | -8.3  | 93    | 0.00     |
| 67 t  | Bromoform                   | 0.117 | 0.122 | -4.3  | 92    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 0.341 | 0.328 | 3.8   | 90    | 0.00     |
| 69 T  | Isopropylbenzene            | 1.268 | 1.345 | -6.1  | 94    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 0.231 | 0.232 | -0.4  | 93    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 0.190 | 0.176 | 7.4   | 90    | 0.00     |
| 72 t  | Bromobenzene                | 0.279 | 0.283 | -1.4  | 91    | 0.00     |
| 73 t  | n-propylbenzene             | 0.319 | 0.341 | -6.9  | 92    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 0.296 | 0.308 | -4.1  | 92    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 1.089 | 1.136 | -4.3  | 93    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 0.297 | 0.315 | -6.1  | 94    | 0.00     |
| 77 t  | tert-Butylbenzene           | 1.022 | 1.049 | -2.6  | 91    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 1.105 | 1.152 | -4.3  | 93    | 0.00     |
| 79 t  | sec-Butylbenzene            | 1.306 | 1.318 | -0.9  | 91    | 0.00     |
| 80    | Nitrobenzene                | 0.005 | 0.005 | 0.0   | 93    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 1.102 | 1.134 | -2.9  | 90    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 0.564 | 0.569 | -0.9  | 92    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 0.551 | 0.563 | -2.2  | 93    | 0.00     |
| 84 t  | n-Butylbenzene              | 1.053 | 1.126 | -6.9  | 92    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 0.512 | 0.523 | -2.1  | 94    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 0.032 | 0.036 | -12.5 | 100   | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 0.306 | 0.326 | -6.5  | 93    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 0.196 | 0.196 | 0.0   | 93    | 0.00     |
| 89 t  | Naphthalene                 | 0.472 | 0.461 | 2.3   | 87    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 0.254 | 0.255 | -0.4  | 90    | 0.00     |

(#) = Out of Range

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