

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\VU101823\  
 Data File : VU055798.D  
 Acq On : 18 Oct 2023 13:34  
 Operator : MD/SY  
 Sample : VSTDICV010  
 Misc : 25.0mL/MSVOA\_U/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 ICVVU101823

Quant Time: Oct 19 03:41:12 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U101823DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Oct 19 03:10:04 2023  
 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     | 100   | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.388 | 0.227 | 41.5#   | 53    | 0.00     |
| 3 t   | Chloromethane               | 0.332 | 0.254 | 23.5    | 71    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 0.340 | 0.267 | 21.5    | 73    | 0.00     |
| 5 T   | Bromomethane                | 0.187 | 0.145 | 22.5    | 70    | 0.00     |
| 6 T   | Chloroethane                | 0.165 | 0.108 | 34.5#   | 59    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.449 | 0.273 | 39.2#   | 56    | -0.02    |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 0.299 | 0.245 | 18.1    | 75    | -0.02    |
| 9 Rt  | 1,1-Dichloroethene          | 0.279 | 0.243 | 12.9    | 79    | -0.02    |
| 10 t  | Iodomethane                 | 0.416 | 0.342 | 17.8    | 71    | 0.00     |
| 11 t  | Allyl Chloride              | 0.363 | 0.361 | 0.6     | 89    | 0.00     |
| 12 t  | Acrylonitrile               | 0.052 | 0.132 | -153.8# | 220#  | 0.00     |
| 13 T  | Acetone                     | 0.112 | 0.057 | 49.1#   | 47    | 0.01     |
| 14 T  | Carbon Disulfide            | 0.910 | 0.717 | 21.2    | 71    | -0.02    |
| 15 RT | Methylene Chloride          | 0.329 | 0.293 | 10.9    | 85    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 0.296 | 0.272 | 8.1     | 82    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 0.550 | 0.515 | 6.4     | 85    | 0.00     |
| 18 T  | 2-Butanone                  | 0.120 | 0.100 | 16.7    | 72    | 0.02     |
| 19    | Cyclohexane                 | 0.469 | 0.424 | 9.6     | 76    | 0.00     |
| 20    | Methylcyclohexane           | 0.487 | 0.474 | 2.7     | 79    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 0.463 | 0.439 | 5.2     | 86    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 0.315 | 0.312 | 1.0     | 86    | 0.00     |
| 23 t  | Diethyl Ether               | 0.167 | 0.139 | 16.8    | 76    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 0.017 | 0.009 | 47.1#   | 47    | 0.12     |
| 25 t  | Methyl tert-Butyl Ether     | 0.600 | 0.607 | -1.2    | 88    | 0.00     |
| 26 t  | Bromochloromethane          | 0.134 | 0.119 | 11.2    | 81    | 0.00     |
| 27 t  | Chloroform                  | 0.564 | 0.527 | 6.6     | 86    | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 0.506 | 0.461 | 8.9     | 82    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 0.400 | 0.385 | 3.8     | 84    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 0.441 | 0.398 | 9.8     | 81    | 0.00     |
| 31 t  | Isopropyl Ether             | 0.739 | 0.780 | -5.5    | 90    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 0.713 | 0.009 | 98.7#   | 1#    | 0.16     |
| 33    | Tert-Amyl methyl ether      | 0.611 | 0.000 | 100.0#  | 0#    | -5.94#   |
| 34 t  | Propionitrile               | 0.019 | 0.036 | -89.5#  | 163   | 0.01     |
| 35 RT | Benzene                     | 1.213 | 1.156 | 4.7     | 85    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 0.329 | 0.305 | 7.3     | 85    | 0.00     |
| 37 RT | Trichloroethene             | 0.316 | 0.288 | 8.9     | 83    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 0.317 | 0.301 | 5.0     | 86    | 0.00     |
| 39 t  | Methacrylonitrile           | 0.071 | 0.070 | 1.4     | 83    | 0.01     |
| 40 t  | Methyl acrylate             | 0.114 | 0.144 | -26.3   | 91    | 0.00     |
| 41 t  | Tetrahydrofuran             | 0.040 | 0.098 | -145.0# | 227#  | 0.00     |
| 42 t  | 1-Chlorobutane              | 0.534 | 0.528 | 1.1     | 86    | 0.00     |
| 43 T  | Dibromomethane              | 0.151 | 0.151 | 0.0     | 93    | 0.00     |
| 44 T  | Bromodichloromethane        | 0.398 | 0.395 | 0.8     | 90    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 0.156 | 0.152 | 2.6     | 84    | 0.03     |
| 46 t  | t-1,4-Dichloro-2-butene     | 0.062 | 0.061 | 1.6     | 95    | 0.00     |
| 47 t  | Methyl methacrylate         | 0.124 | 0.066 | 46.8#   | 44    | 0.00     |
| 48 t  | Ethyl methacrylate          | 0.236 | 0.270 | -14.4   | 94    | 0.00     |
| 49 Rt | Toluene                     | 0.718 | 0.746 | -3.9    | 89    | 0.00     |

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Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 ICVVU101823

Quant Time: Oct 19 03:41:12 2023  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U101823DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Oct 19 03:10:04 2023  
 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       | 0.352 | 0.377 | -7.1  | 94    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 0.431 | 0.452 | -4.9  | 91    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 0.215 | 0.213 | 0.9   | 94    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 0.368 | 0.366 | 0.5   | 92    | 0.00     |
| 54 t  | 2-Hexanone                  | 0.174 | 0.113 | 35.1# | 55    | 0.03     |
| 55 t  | Dibromochloromethane        | 0.260 | 0.254 | 2.3   | 90    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 0.197 | 0.201 | -2.0  | 93    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 0.394 | 0.416 | -5.6  | 106   | 0.00     |
| 58 RT | Tetrachloroethene           | 0.328 | 0.284 | 13.4  | 83    | 0.00     |
| 59 Rt | Chlorobenzene               | 0.793 | 0.785 | 1.0   | 89    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 0.283 | 0.285 | -0.7  | 92    | 0.00     |
| 61 t  | Pentachloroethane           | 0.191 | 0.210 | -9.9  | 98    | 0.00     |
| 62 t  | Hexachloroethane            | 0.202 | 0.222 | -9.9  | 88    | 0.00     |
| 63 Rt | Ethyl Benzene               | 1.346 | 1.474 | -9.5  | 92    | 0.00     |
| 64 RT | m/p-Xylenes                 | 0.517 | 0.569 | -10.1 | 92    | 0.00     |
| 65 RT | o-Xylene                    | 0.508 | 0.544 | -7.1  | 92    | 0.00     |
| 66 RT | Styrene                     | 0.794 | 0.902 | -13.6 | 93    | 0.00     |
| 67 t  | Bromoform                   | 0.139 | 0.143 | -2.9  | 92    | 0.01     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 0.366 | 0.360 | 1.6   | 95    | 0.00     |
| 69 T  | Isopropylbenzene            | 1.307 | 1.456 | -11.4 | 93    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 0.252 | 0.243 | 3.6   | 89    | 0.01     |
| 71 T  | 1,2,3-Trichloropropane      | 0.208 | 0.183 | 12.0  | 81    | 0.00     |
| 72 t  | Bromobenzene                | 0.312 | 0.321 | -2.9  | 92    | 0.00     |
| 73 t  | n-propylbenzene             | 0.337 | 0.382 | -13.4 | 88    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 0.313 | 0.337 | -7.7  | 93    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 1.117 | 1.267 | -13.4 | 91    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 0.319 | 0.335 | -5.0  | 89    | 0.00     |
| 77 t  | tert-Butylbenzene           | 1.024 | 1.144 | -11.7 | 90    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 1.106 | 1.221 | -10.4 | 86    | 0.00     |
| 79 t  | sec-Butylbenzene            | 1.293 | 1.475 | -14.1 | 87    | 0.00     |
| 80    | Nitrobenzene                | 0.007 | 0.002 | 71.4# | 18#   | 0.00     |
| 81 t  | p-Isopropyltoluene          | 1.041 | 1.180 | -13.4 | 84    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 0.636 | 0.635 | 0.2   | 87    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 0.616 | 0.621 | -0.8  | 87    | 0.00     |
| 84 t  | n-Butylbenzene              | 0.888 | 0.958 | -7.9  | 78    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 0.582 | 0.563 | 3.3   | 84    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 0.036 | 0.037 | -2.8  | 82    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 0.316 | 0.347 | -9.8  | 83    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 0.219 | 0.217 | 0.9   | 80    | 0.00     |
| 89 t  | Naphthalene                 | 0.459 | 0.541 | -17.9 | 87    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 0.271 | 0.302 | -11.4 | 85    | 0.00     |

(#) = Out of Range

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