

Data Path : Z:\voasrv\HPCHEM1\MSVOA\_U\Data\U072125\  
 Data File : U063546.D  
 Acq On : 21 Jul 2025 09:13  
 Operator : MD/SY  
 Sample : VSTDCCC010  
 Misc : 25mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Jul 22 04:06:02 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U071625DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Jul 17 03:49:40 2025  
 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    | 99    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.295 | 0.267 | 9.5    | 89    | 0.00     |
| 3 t   | Chloromethane               | 0.274 | 0.261 | 4.7    | 97    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 0.349 | 0.309 | 11.5   | 87    | 0.00     |
| 5 T   | Bromomethane                | 0.263 | 0.239 | 9.1    | 97    | 0.00     |
| 6 T   | Chloroethane                | 0.211 | 0.190 | 10.0   | 88    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.501 | 0.458 | 8.6    | 89    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 0.262 | 0.242 | 7.6    | 90    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 0.260 | 0.229 | 11.9   | 86    | 0.00     |
| 10 t  | Iodomethane                 | 0.314 | 0.243 | 22.6   | 65    | 0.00     |
| 11 t  | Allyl Chloride              | 0.375 | 0.332 | 11.5   | 86    | 0.00     |
| 12 t  | Acrylonitrile               | 0.065 | 0.060 | 7.7    | 95    | 0.00     |
| 13 T  | Acetone                     | 0.052 | 0.075 | -44.2# | 149   | 0.00     |
| 14 T  | Carbon Disulfide            | 0.839 | 0.685 | 18.4   | 79    | 0.00     |
| 15 RT | Methylene Chloride          | 0.306 | 0.266 | 13.1   | 90    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 0.295 | 0.262 | 11.2   | 87    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 0.546 | 0.495 | 9.3    | 90    | 0.00     |
| 18 T  | 2-Butanone                  | 0.081 | 0.093 | -14.8  | 115   | 0.00     |
| 19    | Cyclohexane                 | 0.463 | 0.425 | 8.2    | 91    | 0.00     |
| 20    | Methylcyclohexane           | 0.450 | 0.432 | 4.0    | 87    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 0.465 | 0.427 | 8.2    | 93    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 0.320 | 0.295 | 7.8    | 92    | 0.00     |
| 23 t  | Diethyl Ether               | 0.198 | 0.186 | 6.1    | 90    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 0.024 | 0.022 | 8.3    | 92    | -0.02    |
| 25 t  | Methyl tert-Butyl Ether     | 0.678 | 0.637 | 6.0    | 91    | 0.00     |
| 26 t  | Bromochloromethane          | 0.134 | 0.123 | 8.2    | 88    | 0.00     |
| 27 t  | Chloroform                  | 0.558 | 0.508 | 9.0    | 92    | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 0.473 | 0.471 | 0.4    | 97    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 0.438 | 0.420 | 4.1    | 96    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 0.381 | 0.391 | -2.6   | 99    | 0.00     |
| 31 t  | Isopropyl Ether             | 0.825 | 0.741 | 10.2   | 88    | 0.00     |
| 32    | Ethyl-t-butyl ether         | 0.763 | 0.690 | 9.6    | 89    | 0.00     |
| 33    | Tert-Amyl methyl ether      | 0.663 | 0.614 | 7.4    | 92    | 0.00     |
| 34 t  | Propionitrile               | 0.025 | 0.022 | 12.0   | 95    | 0.00     |
| 35 RT | Benzene                     | 1.178 | 1.237 | -5.0   | 102   | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 0.366 | 0.368 | -0.5   | 97    | 0.00     |
| 37 RT | Trichloroethene             | 0.301 | 0.288 | 4.3    | 89    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 0.276 | 0.296 | -7.2   | 93    | 0.00     |
| 39 t  | Methacrylonitrile           | 0.109 | 0.093 | 14.7   | 91    | 0.00     |
| 40 t  | Methyl acrylate             | 0.159 | 0.141 | 11.3   | 83    | 0.00     |
| 41 t  | Tetrahydrofuran             | 0.057 | 0.048 | 15.8   | 91    | 0.00     |
| 42 t  | 1-Chlorobutane              | 0.583 | 0.567 | 2.7    | 95    | 0.00     |
| 43 T  | Dibromomethane              | 0.143 | 0.149 | -4.2   | 96    | 0.00     |
| 44 T  | Bromodichloromethane        | 0.322 | 0.338 | -5.0   | 102   | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 0.147 | 0.161 | -9.5   | 105   | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 0.055 | 0.068 | -23.6  | 107   | 0.00     |
| 47 t  | Methyl methacrylate         | 0.120 | 0.133 | -10.8  | 104   | 0.00     |
| 48 t  | Ethyl methacrylate          | 0.229 | 0.234 | -2.2   | 92    | 0.00     |
| 49 Rt | Toluene                     | 0.643 | 0.606 | 5.8    | 90    | 0.00     |

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 Misc : 25mL/MSVOA\_U/WATER  
 ALS Vial : 2 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 LabSampleId :  
 VSTDCCC010

Quant Time: Jul 22 04:06:02 2025  
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA\_U\Method\524U071625DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Thu Jul 17 03:49:40 2025  
 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       | 0.253 | 0.247 | 2.4   | 86    | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 0.339 | 0.376 | -10.9 | 103   | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 0.192 | 0.186 | 3.1   | 91    | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 0.332 | 0.321 | 3.3   | 93    | 0.00     |
| 54 t  | 2-Hexanone                  | 0.098 | 0.111 | -13.3 | 107   | 0.00     |
| 55 t  | Dibromochloromethane        | 0.203 | 0.200 | 1.5   | 86    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 0.172 | 0.164 | 4.7   | 91    | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 0.379 | 0.428 | -12.9 | 115   | 0.00     |
| 58 RT | Tetrachloroethene           | 0.275 | 0.243 | 11.6  | 84    | 0.00     |
| 59 Rt | Chlorobenzene               | 0.739 | 0.696 | 5.8   | 90    | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 0.238 | 0.227 | 4.6   | 89    | 0.00     |
| 61 t  | Pentachloroethane           | 0.168 | 0.165 | 1.8   | 88    | 0.00     |
| 62 t  | Hexachloroethane            | 0.166 | 0.175 | -5.4  | 90    | 0.00     |
| 63 Rt | Ethyl Benzene               | 1.271 | 1.201 | 5.5   | 91    | 0.00     |
| 64 RT | m/p-Xylenes                 | 0.494 | 0.478 | 3.2   | 91    | 0.00     |
| 65 RT | o-Xylene                    | 0.475 | 0.467 | 1.7   | 92    | 0.00     |
| 66 RT | Styrene                     | 0.757 | 0.764 | -0.9  | 92    | 0.00     |
| 67 t  | Bromoform                   | 0.104 | 0.103 | 1.0   | 86    | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 0.350 | 0.349 | 0.3   | 97    | 0.00     |
| 69 T  | Isopropylbenzene            | 1.148 | 1.134 | 1.2   | 92    | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 0.234 | 0.239 | -2.1  | 98    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 0.190 | 0.175 | 7.9   | 84    | 0.00     |
| 72 t  | Bromobenzene                | 0.296 | 0.287 | 3.0   | 92    | 0.00     |
| 73 t  | n-propylbenzene             | 0.346 | 0.350 | -1.2  | 93    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 0.305 | 0.301 | 1.3   | 93    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 1.097 | 1.103 | -0.5  | 93    | 0.00     |
| 76 t  | 4-Chlorotoluene             | 0.310 | 0.313 | -1.0  | 94    | 0.00     |
| 77 t  | tert-Butylbenzene           | 1.048 | 1.041 | 0.7   | 92    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 1.088 | 1.105 | -1.6  | 93    | 0.00     |
| 79 t  | sec-Butylbenzene            | 1.413 | 1.430 | -1.2  | 94    | 0.00     |
| 80    | Nitrobenzene                | 0.004 | 0.004 | 0.0   | 70    | 0.00     |
| 81 t  | p-Isopropyltoluene          | 1.180 | 1.201 | -1.8  | 94    | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 0.593 | 0.588 | 0.8   | 93    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 0.593 | 0.593 | 0.0   | 93    | 0.00     |
| 84 t  | n-Butylbenzene              | 1.094 | 1.121 | -2.5  | 92    | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 0.557 | 0.559 | -0.4  | 93    | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 0.033 | 0.036 | -9.1  | 89    | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 0.346 | 0.327 | 5.5   | 84    | 0.00     |
| 88 t  | Hexachlorobutadiene         | 0.197 | 0.187 | 5.1   | 88    | 0.00     |
| 89 t  | Naphthalene                 | 0.621 | 0.545 | 12.2  | 80    | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 0.320 | 0.295 | 7.8   | 83    | 0.00     |

(#) = Out of Range

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