

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU102318\  
 Data File : VU027752.D  
 Acq On : 22 Oct 2018 20:21  
 Operator : MD/SY  
 Sample : VSTDICV010  
 Misc : 25.0mL/MSVOA U/WATER  
 ALS Vial : 9 Sample Multiplier: 1

Instrument :  
 MSVOA\_U  
 ClientSampleId :  
 ICVVU102318

Quant Time: Oct 23 09:17:01 2018  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U0102318DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Oct 23 09:12:15 2018  
 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                    | AvqRF | CCRF  | %Dev    | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|---------|-------|----------|
| 1 i   | Fluorobenzene               | 1.000 | 1.000 | 0.0     | 98    | 0.00     |
| 2 T   | Dichlorodifluoromethane     | 0.397 | 0.246 | 38.0#   | 60    | 0.00     |
| 3 t   | Chloromethane               | 0.425 | 0.349 | 17.9    | 83    | 0.00     |
| 4 Rt  | Vinyl Chloride              | 0.396 | 0.345 | 12.9    | 84    | 0.00     |
| 5 T   | Bromomethane                | 0.208 | 0.192 | 7.7     | 93    | 0.00     |
| 6 T   | Chloroethane                | 0.215 | 0.201 | 6.5     | 93    | 0.00     |
| 7 T   | Trichlorofluoromethane      | 0.483 | 0.450 | 6.8     | 91    | 0.00     |
| 8     | 1,1,2-Trichloro-1,2,2-trifl | 0.254 | 0.232 | 8.7     | 89    | 0.00     |
| 9 Rt  | 1,1-Dichloroethene          | 0.223 | 0.230 | -3.1    | 99    | 0.00     |
| 10 t  | Iodomethane                 | 0.300 | 0.333 | -11.0   | 96    | 0.00     |
| 11 t  | Allyl Chloride              | 0.528 | 0.537 | -1.7    | 101   | 0.00     |
| 12 t  | Acrylonitrile               | 0.068 | 0.160 | -135.3# | 255#  | 0.00     |
| 13 T  | Acetone                     | 0.062 | 0.058 | 6.5     | 102   | 0.00     |
| 14 T  | Carbon Disulfide            | 0.830 | 0.804 | 3.1     | 96    | 0.00     |
| 15 RT | Methylene Chloride          | 0.281 | 0.257 | 8.5     | 98    | 0.00     |
| 16 RT | trans-1,2-Dichloroethene    | 0.248 | 0.254 | -2.4    | 99    | 0.00     |
| 17 t  | 1,1-Dichloroethane          | 0.548 | 0.543 | 0.9     | 99    | 0.00     |
| 18 T  | 2-Butanone                  | 0.103 | 0.099 | 3.9     | 100   | 0.00     |
| 19    | Cyclohexane                 | 0.491 | 0.473 | 3.7     | 95    | 0.00     |
| 20    | Methylcyclohexane           | 0.498 | 0.495 | 0.6     | 95    | 0.00     |
| 21 T  | 2,2-Dichloropropane         | 0.484 | 0.461 | 4.8     | 97    | 0.00     |
| 22 RT | cis-1,2-Dichloroethene      | 0.297 | 0.287 | 3.4     | 96    | 0.00     |
| 23 t  | Diethyl Ether               | 0.212 | 0.203 | 4.2     | 93    | 0.00     |
| 24 t  | tert-Butyl Alcohol          | 0.023 | 0.011 | 52.2#   | 46    | 0.01     |
| 25 t  | Methyl tert-Butyl Ether     | 0.710 | 0.705 | 0.7     | 97    | 0.00     |
| 26 t  | Bromochloromethane          | 0.119 | 0.113 | 5.0     | 93    | 0.00     |
| 27 t  | Chloroform                  | 0.533 | 0.542 | -1.7    | 100   | 0.00     |
| 28 RT | 1,1,1-Trichloroethane       | 0.475 | 0.483 | -1.7    | 98    | 0.00     |
| 29 T  | 1,1-Dichloropropene         | 0.415 | 0.417 | -0.5    | 96    | 0.00     |
| 30 RT | Carbon Tetrachloride        | 0.424 | 0.430 | -1.4    | 99    | 0.00     |
| 31 t  | Isopropyl Ether             | 0.987 | 1.065 | -7.9    | 104   | 0.00     |
| 32    | Ethyl-t-butyl ether         | 0.876 | 0.000 | 100.0#  | 0#    | -4.08#   |
| 33    | Tert-Amyl methyl ether      | 0.743 | 0.000 | 100.0#  | 0#    | -5.59#   |
| 34 t  | Propionitrile               | 0.025 | 0.054 | -116.0# | 209#  | 0.00     |
| 35 RT | Benzene                     | 1.159 | 1.183 | -2.1    | 98    | 0.00     |
| 36 RT | 1,2-Dichloroethane          | 0.386 | 0.385 | 0.3     | 97    | 0.00     |
| 37 RT | Trichloroethene             | 0.286 | 0.285 | 0.3     | 98    | 0.00     |
| 38 Rt | 1,2-Dichloropropane         | 0.325 | 0.325 | 0.0     | 96    | 0.00     |
| 39 t  | Methacrylonitrile           | 0.136 | 0.137 | -0.7    | 99    | 0.00     |
| 40 t  | Methyl acrylate             | 0.177 | 0.191 | -7.9    | 103   | 0.00     |
| 41 t  | Tetrahydrofuran             | 0.075 | 0.166 | -121.3# | 236#  | 0.00     |
| 42 t  | 1-Chlorobutane              | 0.654 | 0.655 | -0.2    | 97    | 0.00     |
| 43 T  | Dibromomethane              | 0.148 | 0.156 | -5.4    | 99    | 0.00     |
| 44 T  | Bromodichloromethane        | 0.402 | 0.411 | -2.2    | 99    | 0.00     |
| 45 T  | 4-Methyl-2-Pentanone        | 0.240 | 0.246 | -2.5    | 99    | 0.00     |
| 46 t  | t-1,4-Dichloro-2-butene     | 0.087 | 0.097 | -11.5   | 105   | -0.02    |

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 Misc : 25.0mL/MSVOA U/WATER  
 ALS Vial : 9 Sample Multiplier: 1

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 MSVOA\_U  
 ClientSampleId :  
 ICVVU102318

Quant Time: Oct 23 09:17:01 2018  
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U0102318DW.M  
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER  
 QLast Update : Tue Oct 23 09:12:15 2018  
 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                    | AvqRF | CCRF  | %Dev  | Area% | Dev(min) |
|-------|-----------------------------|-------|-------|-------|-------|----------|
| 47 t  | Methyl methacrylate         | 0.156 | 0.084 | 46.2# | 50    | 0.00     |
| 48 t  | Ethyl methacrylate          | 0.324 | 0.334 | -3.1  | 96    | 0.00     |
| 49 Rt | Toluene                     | 0.688 | 0.732 | -6.4  | 101   | 0.00     |
| 50 T  | t-1,3-Dichloropropene       | 0.400 | 0.429 | -7.2  | 101   | 0.00     |
| 51 T  | cis-1,3-Dichloropropene     | 0.472 | 0.492 | -4.2  | 99    | 0.00     |
| 52 RT | 1,1,2-Trichloroethane       | 0.200 | 0.214 | -7.0  | 102   | 0.00     |
| 53 t  | 1,3-Dichloropropane         | 0.384 | 0.394 | -2.6  | 99    | 0.00     |
| 54 t  | 2-Hexanone                  | 0.171 | 0.179 | -4.7  | 101   | 0.00     |
| 55 t  | Dibromochloromethane        | 0.251 | 0.258 | -2.8  | 97    | 0.00     |
| 56 T  | 1,2-Dibromoethane           | 0.196 | 0.205 | -4.6  | 100   | 0.00     |
| 57 S  | 4-Bromofluorobenzene        | 0.379 | 0.377 | 0.5   | 95    | 0.00     |
| 58 RT | Tetrachloroethene           | 0.270 | 0.273 | -1.1  | 100   | 0.00     |
| 59 Rt | Chlorobenzene               | 0.737 | 0.768 | -4.2  | 100   | 0.00     |
| 60 T  | 1,1,1,2-Tetrachloroethane   | 0.268 | 0.280 | -4.5  | 99    | 0.00     |
| 61 t  | Pentachloroethane           | 0.200 | 0.223 | -11.5 | 101   | 0.00     |
| 62 t  | Hexachloroethane            | 0.220 | 0.243 | -10.5 | 101   | 0.00     |
| 63 Rt | Ethyl Benzene               | 1.396 | 1.474 | -5.6  | 100   | 0.00     |
| 64 RT | m/p-Xylenes                 | 0.498 | 0.530 | -6.4  | 101   | 0.00     |
| 65 RT | o-Xylene                    | 0.494 | 0.514 | -4.0  | 99    | 0.00     |
| 66 RT | Styrene                     | 0.815 | 0.850 | -4.3  | 97    | 0.00     |
| 67 t  | Bromoform                   | 0.145 | 0.159 | -9.7  | 101   | 0.00     |
| 68 S  | 1,2-Dichlorobenzene-d4      | 0.358 | 0.357 | 0.3   | 97    | 0.00     |
| 69 T  | Isopropylbenzene            | 1.352 | 1.423 | -5.3  | 101   | 0.00     |
| 70 T  | 1,1,2,2-Tetrachloroethane   | 0.266 | 0.273 | -2.6  | 99    | 0.00     |
| 71 T  | 1,2,3-Trichloropropane      | 0.206 | 0.137 | 33.5# | 64    | 0.00     |
| 72 t  | Bromobenzene                | 0.301 | 0.327 | -8.6  | 100   | 0.00     |
| 73 t  | n-propylbenzene             | 0.349 | 0.371 | -6.3  | 98    | 0.00     |
| 74 t  | 2-Chlorotoluene             | 0.299 | 0.315 | -5.4  | 99    | 0.00     |
| 75 t  | 1,3,5-Trimethylbenzene      | 1.128 | 1.213 | -7.5  | 102   | 0.00     |
| 76 t  | 4-Chlorotoluene             | 0.303 | 0.328 | -8.3  | 102   | 0.00     |
| 77 t  | tert-Butylbenzene           | 1.110 | 1.152 | -3.8  | 99    | 0.00     |
| 78 t  | 1,2,4-Trimethylbenzene      | 1.159 | 1.250 | -7.9  | 101   | 0.00     |
| 79 t  | sec-Butylbenzene            | 1.470 | 1.491 | -1.4  | 96    | 0.00     |
| 80    | Nitrobenzene                | 0.010 | 0.002 | 80.0# | 19#   | 0.00     |
| 81 t  | p-Isopropyltoluene          | 1.201 | 1.302 | -8.4  | 101   | 0.00     |
| 82 t  | 1,3-Dichlorobenzene         | 0.596 | 0.631 | -5.9  | 99    | 0.00     |
| 83 Rt | 1,4-Dichlorobenzene         | 0.597 | 0.641 | -7.4  | 104   | 0.00     |
| 84 t  | n-Butylbenzene              | 1.167 | 1.307 | -12.0 | 101   | 0.00     |
| 85 Rt | 1,2-Dichlorobenzene         | 0.569 | 0.595 | -4.6  | 100   | 0.00     |
| 86 t  | 1,2-Dibromo-3-Chloropropane | 0.049 | 0.054 | -10.2 | 100   | 0.00     |
| 87 Rt | 1,2,4-Trichlorobenzene      | 0.413 | 0.468 | -13.3 | 101   | 0.00     |
| 88 t  | Hexachlorobutadiene         | 0.248 | 0.265 | -6.9  | 101   | 0.00     |
| 89 t  | Naphthalene                 | 0.776 | 0.886 | -14.2 | 104   | 0.00     |
| 90 t  | 1,2,3-Trichlorobenzene      | 0.393 | 0.434 | -10.4 | 104   | 0.00     |

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Max. RRF Dev : 30% Max. Rel. Area : 200%

| Compound           | AvgRF                        | CCRF | %Dev Area | % Dev(min) |
|--------------------|------------------------------|------|-----------|------------|
| -----              |                              |      |           |            |
| (#) = Out of Range | SPCC's out = 0 CCC's out = 0 |      |           |            |