

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\U081722\
 Data File : VU050366.D
 Acq On : 17 Aug 2022 20:47
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampled :
 VSTDCCC010

Quant Time: Aug 18 06:11:26 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081722DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Aug 18 06:06:02 2022
 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 | 94 | 0.00 |
| 2 T | Dichlorodifluoromethane | 10.000 | 7.478 | 25.2 | 75 | 0.00 |
| 3 t | Chloromethane | 10.000 | 9.053 | 9.5 | 92 | 0.00 |
| 4 Rt | Vinyl Chloride | 10.000 | 9.122 | 8.8 | 89 | 0.00 |
| 5 T | Bromomethane | 10.000 | 10.040 | -0.4 | 100 | 0.00 |
| 6 T | Chloroethane | 10.000 | 9.685 | 3.1 | 96 | 0.00 |
| 7 T | Trichlorofluoromethane | 10.000 | 8.360 | 16.4 | 83 | 0.00 |
| 8 | 1,1,2-Trichloro-1,2,2-trifl | 10.000 | 7.617 | 23.8 | 76 | 0.00 |
| 9 Rt | 1,1-Dichloroethene | 10.000 | 9.006 | 9.9 | 88 | 0.00 |
| 10 t | Iodomethane | 10.000 | 11.974 | -19.7 | 102 | 0.00 |
| 11 t | Allyl Chloride | 10.000 | 9.613 | 3.9 | 96 | 0.00 |
| 12 t | Acrylonitrile | 20.000 | 19.132 | 4.3 | 93 | 0.00 |
| 13 T | Acetone | 50.000 | 54.836 | -9.7 | 99 | 0.00 |
| 14 T | Carbon Disulfide | 10.000 | 8.849 | 11.5 | 89 | 0.00 |
| 15 RT | Methylene Chloride | 10.000 | 6.969 | 30.3# | 109 | 0.00 |
| 16 RT | trans-1,2-Dichloroethene | 10.000 | 9.672 | 3.3 | 93 | -0.02 |
| 17 t | 1,1-Dichloroethane | 10.000 | 10.410 | -4.1 | 98 | -0.03 |
| 18 T | 2-Butanone | 50.000 | 54.897 | -9.8 | 101 | 0.00 |
| 19 | Cyclohexane | 10.000 | 8.157 | 18.4 | 73 | 0.00 |
| 20 | Methylcyclohexane | 10.000 | 8.113 | 18.9 | 71 | -0.03 |
| 21 T | 2,2-Dichloropropane | 10.000 | 8.981 | 10.2 | 82 | 0.00 |
| 22 RT | cis-1,2-Dichloroethene | 10.000 | 10.860 | -8.6 | 103 | 0.00 |
| 23 t | Diethyl Ether | 10.000 | 9.519 | 4.8 | 97 | 0.00 |
| 24 t | tert-Butyl Alcohol | 100.000 | 113.176 | -13.2 | 107 | 0.03 |
| 25 t | Methyl tert-Butyl Ether | 10.000 | 10.175 | -1.8 | 97 | -0.01 |
| 26 t | Bromochloromethane | 10.000 | 10.222 | -2.2 | 108 | 0.00 |
| 27 t | Chloroform | 10.000 | 10.409 | -4.1 | 104 | 0.00 |
| 28 RT | 1,1,1-Trichloroethane | 10.000 | 9.653 | 3.5 | 93 | 0.00 |
| 29 T | 1,1-Dichloropropene | 10.000 | 10.065 | -0.6 | 90 | 0.00 |
| 30 RT | Carbon Tetrachloride | 10.000 | 9.385 | 6.2 | 91 | 0.00 |
| 31 t | Isopropyl Ether | 10.000 | 11.031 | -10.3 | 99 | -0.02 |
| 32 | Ethyl-t-butyl ether | 10.000 | 11.617 | -16.2 | 98 | -0.01 |
| 33 | Tert-Amyl methyl ether | 10.000 | 11.978 | -19.8 | 103 | 0.00 |
| 34 t | Propionitrile | 50.000 | 53.741 | -7.5 | 97 | 0.00 |
| 35 RT | Benzene | 10.000 | 10.414 | -4.1 | 96 | 0.00 |
| 36 RT | 1,2-Dichloroethane | 10.000 | 10.110 | -1.1 | 97 | 0.00 |
| 37 RT | Trichloroethene | 10.000 | 9.895 | 1.1 | 94 | -0.03 |
| 38 Rt | 1,2-Dichloropropane | 10.000 | 10.570 | -5.7 | 98 | -0.03 |
| 39 t | Methacrylonitrile | 10.000 | 12.173 | -21.7 | 99 | 0.00 |
| 40 t | Methyl acrylate | 10.000 | 12.219 | -22.2 | 101 | 0.00 |
| 41 t | Tetrahydrofuran | 20.000 | 22.933 | -14.7 | 99 | 0.00 |
| 42 t | 1-Chlorobutane | 10.000 | 9.912 | 0.9 | 91 | 0.00 |
| 43 T | Dibromomethane | 10.000 | 10.122 | -1.2 | 97 | -0.03 |
| 44 T | Bromodichloromethane | 10.000 | 10.388 | -3.9 | 98 | -0.02 |
| 45 T | 4-Methyl-2-Pentanone | 50.000 | 54.173 | -8.3 | 96 | -0.01 |
| 46 t | t-1,4-Dichloro-2-butene | 20.000 | 20.730 | -3.7 | 79 | 0.00 |
| 47 t | Methyl methacrylate | 20.000 | 22.617 | -13.1 | 93 | -0.02 |
| 48 t | Ethyl methacrylate | 10.000 | 9.751 | 2.5 | 92 | 0.00 |
| 49 Rt | Toluene | 10.000 | 11.344 | -13.4 | 97 | -0.02 |

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 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Aug 18 06:11:26 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081722DW.M
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 QLast Update : Thu Aug 18 06:06:02 2022
 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 | 10.578 | -5.8 | 93 | -0.02 |
| 51 T cis-1,3-Dichloropropene | 10.000 | 10.440 | -4.4 | 93 | -0.02 |
| 52 RT 1,1,2-Trichloroethane | 10.000 | 10.020 | -0.2 | 98 | 0.00 |
| 53 t 1,3-Dichloropropane | 10.000 | 10.598 | -6.0 | 97 | 0.00 |
| 54 t 2-Hexanone | 50.000 | 55.292 | -10.6 | 94 | 0.00 |
| 55 t Dibromochloromethane | 10.000 | 10.349 | -3.5 | 98 | 0.00 |
| 56 T 1,2-Dibromoethane | 10.000 | 10.597 | -6.0 | 97 | 0.00 |
| 57 S 4-Bromofluorobenzene | 1.000 | 1.090 | -9.0 | 97 | 0.00 |
| 58 RT Tetrachloroethene | 10.000 | 9.400 | 6.0 | 91 | -0.01 |
| 59 Rt Chlorobenzene | 10.000 | 10.742 | -7.4 | 96 | 0.00 |
| 60 T 1,1,1,2-Tetrachloroethane | 10.000 | 10.445 | -4.5 | 100 | 0.00 |
| 61 t Pentachloroethane | 10.000 | 10.497 | -5.0 | 98 | 0.00 |
| 62 t Hexachloroethane | 10.000 | 10.221 | -2.2 | 96 | 0.00 |
| 63 Rt Ethyl Benzene | 10.000 | 9.764 | 2.4 | 93 | 0.00 |
| 64 RT m/p-Xylenes | 20.000 | 20.176 | -0.9 | 94 | 0.00 |
| 65 RT o-Xylene | 10.000 | 10.179 | -1.8 | 95 | 0.00 |
| 66 RT Styrene | 10.000 | 10.404 | -4.0 | 98 | 0.00 |
| 67 t Bromoform | 10.000 | 10.261 | -2.6 | 94 | 0.00 |
| 68 S 1,2-Dichlorobenzene-d4 | 1.000 | 1.008 | -0.8 | 94 | 0.00 |
| 69 T Isopropylbenzene | 10.000 | 9.672 | 3.3 | 92 | 0.00 |
| 70 T 1,1,2,2-Tetrachloroethane | 10.000 | 9.790 | 2.1 | 92 | 0.00 |
| 71 T 1,2,3-Trichloropropane | 10.000 | 9.667 | 3.3 | 94 | 0.00 |
| 72 t Bromobenzene | 10.000 | 11.037 | -10.4 | 96 | 0.00 |
| 73 t n-propylbenzene | 10.000 | 9.532 | 4.7 | 89 | 0.00 |
| 74 t 2-Chlorotoluene | 10.000 | 10.306 | -3.1 | 97 | 0.00 |
| 75 t 1,3,5-Trimethylbenzene | 10.000 | 10.076 | -0.8 | 95 | 0.00 |
| 76 t 4-Chlorotoluene | 10.000 | 11.948 | -19.5 | 95 | 0.00 |
| 77 t tert-Butylbenzene | 10.000 | 9.637 | 3.6 | 92 | 0.00 |
| 78 t 1,2,4-Trimethylbenzene | 10.000 | 10.191 | -1.9 | 93 | 0.00 |
| 79 t sec-Butylbenzene | 10.000 | 9.096 | 9.0 | 86 | 0.00 |
| 80 Nitrobenzene | 50.000 | 42.709 | 14.6 | 82 | 0.00 |
| 81 t p-Isopropyltoluene | 10.000 | 6.706 | 32.9# | 87 | 0.00 |
| 82 t 1,3-Dichlorobenzene | 10.000 | 10.659 | -6.6 | 93 | 0.00 |
| 83 Rt 1,4-Dichlorobenzene | 10.000 | 10.294 | -2.9 | 92 | 0.00 |
| 84 t n-Butylbenzene | 10.000 | 8.403 | 16.0 | 81 | 0.00 |
| 85 Rt 1,2-Dichlorobenzene | 10.000 | 10.567 | -5.7 | 93 | 0.00 |
| 86 t 1,2-Dibromo-3-Chloropropane | 10.000 | 10.013 | -0.1 | 93 | 0.00 |
| 87 Rt 1,2,4-Trichlorobenzene | 10.000 | 8.623 | 13.8 | 83 | 0.00 |
| 88 t Hexachlorobutadiene | 10.000 | 8.580 | 14.2 | 84 | 0.00 |
| 89 t Naphthalene | 10.000 | 13.121 | -31.2# | 131 | 0.00 |
| 90 t 1,2,3-Trichlorobenzene | 10.000 | 8.966 | 10.3 | 84 | 0.00 |

(#) = Out of Range

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