

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU081022\
 Data File : VU050223.D
 Acq On : 10 Aug 2022 17:10
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Aug 10 18:15:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081022DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Aug 10 15:54:10 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	93	0.00
2 T	Dichlorodifluoromethane	10.000	9.159	8.4	88	0.00
3 t	Chloromethane	10.000	9.041	9.6	96	0.00
4 Rt	Vinyl Chloride	10.000	9.281	7.2	91	0.00
5 T	Bromomethane	10.000	10.629	-6.3	107	0.00
6 T	Chloroethane	10.000	9.293	7.1	95	0.00
7 T	Trichlorofluoromethane	10.000	9.275	7.2	91	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.321	6.8	90	0.00
9 Rt	1,1-Dichloroethene	10.000	9.041	9.6	89	0.00
10 t	Iodomethane	10.000	9.673	3.3	89	0.00
11 t	Allyl Chloride	10.000	9.374	6.3	90	0.00
12 t	Acrylonitrile	20.000	20.271	-1.4	97	0.01
13 T	Acetone	50.000	47.787	4.4	96	0.03
14 T	Carbon Disulfide	10.000	8.537	14.6	90	0.00
15 RT	Methylene Chloride	10.000	8.541	14.6	100	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.633	3.7	93	0.00
17 t	1,1-Dichloroethane	10.000	9.550	4.5	94	0.00
18 T	2-Butanone	50.000	50.243	-0.5	95	0.00
19	Cyclohexane	10.000	8.523	14.8	73	0.00
20	Methylcyclohexane	10.000	10.136	-1.4	86	0.00
21 T	2,2-Dichloropropane	10.000	9.653	3.5	101	0.00
22 RT	cis-1,2-Dichloroethene	10.000	9.818	1.8	95	0.00
23 t	Diethyl Ether	10.000	9.592	4.1	99	0.00
24 t	tert-Butyl Alcohol	100.000	136.107	-36.1#	132	0.04
25 t	Methyl tert-Butyl Ether	10.000	9.979	0.2	95	0.00
26 t	Bromochloromethane	10.000	9.884	1.2	98	0.00
27 t	Chloroform	10.000	9.681	3.2	97	0.00
28 RT	1,1,1-Trichloroethane	10.000	9.564	4.4	93	0.00
29 T	1,1-Dichloropropene	10.000	9.874	1.3	91	0.00
30 RT	Carbon Tetrachloride	10.000	9.443	5.6	92	0.00
31 t	Isopropyl Ether	10.000	10.281	-2.8	93	0.00
32	Ethyl-t-butyl ether	10.000	9.957	0.4	93	0.00
33	Tert-Amyl methyl ether	10.000	10.261	-2.6	94	0.00
34 t	Propionitrile	50.000	50.070	-0.1	94	0.01
35 RT	Benzene	10.000	9.886	1.1	94	0.00
36 RT	1,2-Dichloroethane	10.000	9.699	3.0	97	0.00
37 RT	Trichloroethene	10.000	9.416	5.8	91	0.00
38 Rt	1,2-Dichloropropane	10.000	9.967	0.3	98	0.00
39 t	Methacrylonitrile	10.000	10.641	-6.4	95	0.00
40 t	Methyl acrylate	10.000	10.583	-5.8	92	0.00
41 t	Tetrahydrofuran	20.000	19.090	4.6	94	0.00
42 t	1-Chlorobutane	10.000	10.012	-0.1	91	0.00
43 T	Dibromomethane	10.000	9.742	2.6	96	0.00
44 T	Bromodichloromethane	10.000	9.945	0.5	97	0.00
45 T	4-Methyl-2-Pentanone	50.000	51.337	-2.7	95	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	20.320	-1.6	108	0.00
47 t	Methyl methacrylate	20.000	21.407	-7.0	95	0.00
48 t	Ethyl methacrylate	10.000	10.618	-6.2	93	0.00
49 Rt	Toluene	10.000	10.473	-4.7	95	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU081022\
 Data File : VU050223.D
 Acq On : 10 Aug 2022 17:10
 Operator : SY/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Aug 10 18:15:27 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U081022DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Aug 10 15:54:10 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	10.109	-1.1	95	0.00
51 T	cis-1,3-Dichloropropene	10.000	9.938	0.6	95	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.782	2.2	98	0.00
53 t	1,3-Dichloropropane	10.000	9.943	0.6	98	0.00
54 t	2-Hexanone	50.000	50.634	-1.3	93	0.00
55 t	Dibromochloromethane	10.000	9.855	1.4	98	0.00
56 T	1,2-Dibromoethane	10.000	10.036	-0.4	97	0.00
57 S	4-Bromofluorobenzene	1.000	1.038	-3.8	95	0.00
58 RT	Tetrachloroethene	10.000	9.548	4.5	93	0.00
59 Rt	Chlorobenzene	10.000	10.022	-0.2	93	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	9.772	2.3	96	0.00
61 t	Pentachloroethane	10.000	9.827	1.7	97	0.00
62 t	Hexachloroethane	10.000	9.825	1.8	89	0.00
63 Rt	Ethyl Benzene	10.000	10.701	-7.0	91	0.00
64 RT	m/p-Xylenes	20.000	22.471	-12.4	93	0.00
65 RT	o-Xylene	10.000	10.998	-10.0	94	0.00
66 RT	Styrene	10.000	11.733	-17.3	95	0.00
67 t	Bromoform	10.000	9.699	3.0	95	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.020	-2.0	95	0.00
69 T	Isopropylbenzene	10.000	10.706	-7.1	90	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.699	3.0	97	0.00
71 T	1,2,3-Trichloropropane	10.000	10.901	-9.0	113	0.00
72 t	Bromobenzene	10.000	10.231	-2.3	95	0.00
73 t	n-propylbenzene	10.000	11.266	-12.7	90	0.00
74 t	2-Chlorotoluene	10.000	10.706	-7.1	93	0.00
75 t	1,3,5-Trimethylbenzene	10.000	11.272	-12.7	92	0.00
76 t	4-Chlorotoluene	10.000	11.301	-13.0	95	0.00
77 t	tert-Butylbenzene	10.000	10.435	-4.4	90	0.00
78 t	1,2,4-Trimethylbenzene	10.000	11.217	-12.2	92	0.00
79 t	sec-Butylbenzene	10.000	10.624	-6.2	87	0.00
80	Nitrobenzene	50.000	49.146	1.7	89	0.00
81 t	p-Isopropyltoluene	10.000	10.943	-9.4	88	0.00
82 t	1,3-Dichlorobenzene	10.000	10.086	-0.9	94	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.313	-3.1	95	0.00
84 t	n-Butylbenzene	10.000	10.308	-3.1	87	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.246	-2.5	95	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	9.870	1.3	95	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	9.808	1.9	87	0.00
88 t	Hexachlorobutadiene	10.000	9.408	5.9	86	0.00
89 t	Naphthalene	10.000	10.016	-0.2	85	0.00
90 t	1,2,3-Trichlorobenzene	10.000	9.965	0.4	88	0.00

(#) = Out of Range

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