

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041923\
 Data File : VU053783.D
 Acq On : 19 Apr 2023 15:05
 Operator : JC/MD
 Sample : VSTDICV010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICVVU041923

Quant Time: Apr 21 09:19:09 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U041923DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Apr 21 09:18:21 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 i	Fluorobenzene	1.000	1.000	0.0	104	0.00
2 T	Dichlorodifluoromethane	10.000	8.807	11.9	89	0.00
3 t	Chloromethane	10.000	11.382	-13.8	124	0.00
4 Rt	Vinyl Chloride	10.000	11.723	-17.2	119	0.00
5 T	Bromomethane	10.000	10.868	-8.7	128	0.00
6 T	Chloroethane	10.000	11.313	-13.1	117	0.00
7 T	Trichlorofluoromethane	10.000	10.474	-4.7	108	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.772	2.3	98	0.00
9 Rt	1,1-Dichloroethene	10.000	11.158	-11.6	115	0.00
10 t	Iodomethane	10.000	12.665	-26.6	116	0.00
11 t	Allyl Chloride	10.000	11.776	-17.8	124	0.00
12 t	Acrylonitrile	20.000	46.949	-134.7#	244	0.00
13 T	Acetone	50.000	32.473	35.1#	58	0.00
14 T	Carbon Disulfide	10.000	16.880	-68.8#	175	0.00
15 RT	Methylene Chloride	10.000	8.790	12.1	98	0.00
16 RT	trans-1,2-Dichloroethene	10.000	11.153	-11.5	115	0.00
17 t	1,1-Dichloroethane	10.000	10.243	-2.4	106	0.00
18 T	2-Butanone	50.000	33.411	33.2#	63	0.00
19	Cyclohexane	10.000	11.746	-17.5	123	0.00
20	Methylcyclohexane	10.000	12.611	-26.1	123	0.00
21 T	2,2-Dichloropropane	10.000	10.029	-0.3	103	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.617	-6.2	112	0.00
23 t	Diethyl Ether	10.000	9.817	1.8	98	0.00
24 t	tert-Butyl Alcohol	100.000	42.006	58.0#	41	0.00
25 t	Methyl tert-Butyl Ether	10.000	9.871	1.3	101	0.00
26 t	Bromochloromethane	10.000	7.882	21.2	82	0.00
27 t	Chloroform	10.000	9.847	1.5	102	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.037	-0.4	101	0.00
29 T	1,1-Dichloropropene	10.000	11.167	-11.7	115	0.00
30 RT	Carbon Tetrachloride	10.000	10.443	-4.4	106	0.00
31 t	Isopropyl Ether	10.000	9.920	0.8	103	0.00
32	Ethyl-t-butyl ether	10.000	0.128	98.7#	1	0.15
33	Tert-Amyl methyl ether	10.000	0.289	97.1#	3	-0.18
34 t	Propionitrile	50.000	89.084	-78.2#	204	0.00
35 RT	Benzene	10.000	10.838	-8.4	109	0.00
36 RT	1,2-Dichloroethane	10.000	9.899	1.0	102	0.00
37 RT	Trichloroethene	10.000	10.464	-4.6	104	0.00
38 Rt	1,2-Dichloropropane	10.000	10.115	-1.2	102	0.00
39 t	Methacrylonitrile	10.000	8.125	18.8	95	0.00
40 t	Methyl acrylate	10.000	9.047	9.5	99	0.00
41 t	Tetrahydrofuran	20.000	42.871	-114.4#	253	0.00
42 t	1-Chlorobutane	10.000	11.300	-13.0	116	0.00
43 T	Dibromomethane	10.000	10.566	-5.7	108	0.00
44 T	Bromodichloromethane	10.000	10.349	-3.5	103	0.00
45 T	4-Methyl-2-Pentanone	50.000	47.254	5.5	91	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	21.117	-5.6	91	0.00
47 t	Methyl methacrylate	20.000	9.748	51.3#	48	0.00
48 t	Ethyl methacrylate	10.000	10.282	-2.8	100	0.00
49 Rt	Toluene	10.000	11.038	-10.4	108	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU041923\
 Data File : VU053783.D
 Acq On : 19 Apr 2023 15:05
 Operator : JC/MD
 Sample : VSTDICV010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICSVU041923

Quant Time: Apr 21 09:19:09 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U041923DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Fri Apr 21 09:18:21 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	Amount	Calc.	%Dev	Area%	Dev(min)
50 T	t-1,3-Dichloropropene	10.000	11.009	-10.1	105	0.00
51 T	cis-1,3-Dichloropropene	10.000	10.897	-9.0	106	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.598	4.0	98	0.00
53 t	1,3-Dichloropropane	10.000	10.252	-2.5	103	0.00
54 t	2-Hexanone	50.000	38.670	22.7	67	0.00
55 t	Dibromochloromethane	10.000	10.304	-3.0	99	0.00
56 T	1,2-Dibromoethane	10.000	10.708	-7.1	104	0.00
57 S	4-Bromofluorobenzene	1.000	1.084	-8.4	110	0.00
58 RT	Tetrachloroethene	10.000	10.680	-6.8	106	0.00
59 Rt	Chlorobenzene	10.000	9.987	0.1	98	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.059	-0.6	98	0.00
61 t	Pentachloroethane	10.000	10.480	-4.8	104	0.00
62 t	Hexachloroethane	10.000	10.913	-9.1	101	0.00
63 Rt	Ethyl Benzene	10.000	11.218	-12.2	105	0.00
64 RT	m/p-Xylenes	20.000	23.380	-16.9	108	0.00
65 RT	o-Xylene	10.000	11.082	-10.8	103	0.00
66 RT	Styrene	10.000	11.581	-15.8	104	0.00
67 t	Bromoform	10.000	10.742	-7.4	101	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.074	-7.4	106	0.00
69 T	Isopropylbenzene	10.000	11.251	-12.5	104	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.622	3.8	98	0.00
71 T	1,2,3-Trichloropropane	10.000	8.164	18.4	90	0.00
72 t	Bromobenzene	10.000	10.195	-2.0	99	0.00
73 t	n-propylbenzene	10.000	11.626	-16.3	102	0.00
74 t	2-Chlorotoluene	10.000	10.851	-8.5	104	0.00
75 t	1,3,5-Trimethylbenzene	10.000	11.610	-16.1	105	0.00
76 t	4-Chlorotoluene	10.000	10.995	-9.9	99	0.00
77 t	tert-Butylbenzene	10.000	10.956	-9.6	100	0.00
78 t	1,2,4-Trimethylbenzene	10.000	11.430	-14.3	100	0.00
79 t	sec-Butylbenzene	10.000	11.264	-12.6	100	0.00
80	Nitrobenzene	50.000	9.897	80.2#	19	0.00
81 t	p-Isopropyltoluene	10.000	11.592	-15.9	101	0.00
82 t	1,3-Dichlorobenzene	10.000	10.347	-3.5	98	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.807	-8.1	100	0.00
84 t	n-Butylbenzene	10.000	9.644	3.6	98	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.119	-1.2	95	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	10.019	-0.2	91	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	10.201	-2.0	96	0.00
88 t	Hexachlorobutadiene	10.000	10.067	-0.7	93	0.00
89 t	Naphthalene	10.000	10.268	-2.7	93	0.00
90 t	1,2,3-Trichlorobenzene	10.000	9.679	3.2	89	0.00

(#) = Out of Range

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