

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU021422\
 Data File : VU047131.D
 Acq On : 14 Feb 2022 13:44
 Operator : SY/MD
 Sample : VSTDICV010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICVVU021422

Quant Time: Feb 15 03:18:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U021422DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Feb 15 03:17:28 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	100	0.00
2 T	Dichlorodifluoromethane	10.000	5.858	41.4#	55	0.00
3 t	Chloromethane	10.000	8.086	19.1	76	0.00
4 Rt	Vinyl Chloride	10.000	8.284	17.2	78	0.00
5 T	Bromomethane	10.000	8.644	13.6	87	0.00
6 T	Chloroethane	10.000	9.298	7.0	88	0.00
7 T	Trichlorofluoromethane	10.000	9.432	5.7	88	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	10.282	-2.8	96	0.00
9 Rt	1,1-Dichloroethene	10.000	9.634	3.7	93	0.00
10 t	Iodomethane	10.000	8.556	14.4	86	0.00
11 t	Allyl Chloride	10.000	10.311	-3.1	97	0.00
12 t	Acrylonitrile	20.000	48.272	-141.4#	267	0.00
13 T	Acetone	50.000	79.860	-59.7#	210	0.00
14 T	Carbon Disulfide	10.000	8.541	14.6	80	0.00
15 RT	Methylene Chloride	10.000	10.043	-0.4	97	0.00
16 RT	trans-1,2-Dichloroethene	10.000	10.043	-0.4	94	0.00
17 t	1,1-Dichloroethane	10.000	10.453	-4.5	98	0.00
18 T	2-Butanone	50.000	63.511	-27.0	126	0.00
19	Cyclohexane	10.000	10.173	-1.7	95	0.00
20	Methylcyclohexane	10.000	10.228	-2.3	95	0.00
21 T	2,2-Dichloropropane	10.000	10.484	-4.8	97	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.470	-4.7	100	0.00
23 t	Diethyl Ether	10.000	9.793	2.1	92	0.00
24 t	tert-Butyl Alcohol	100.000	67.453	32.5#	64	0.02
25 t	Methyl tert-Butyl Ether	10.000	10.683	-6.8	102	0.00
26 t	Bromochloromethane	10.000	9.486	5.1	88	0.00
27 t	Chloroform	10.000	10.343	-3.4	100	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.337	-3.4	97	0.00
29 T	1,1-Dichloropropene	10.000	10.070	-0.7	94	0.00
30 RT	Carbon Tetrachloride	10.000	10.404	-4.0	96	0.00
31 t	Isopropyl Ether	10.000	11.038	-10.4	103	0.00
32	Ethyl-t-butyl ether	10.000	0.000	100.0#	0	-4.51#
33	Tert-Amyl methyl ether	10.000	0.000	100.0#	0	-5.94#
34 t	Propionitrile	50.000	87.501	-75.0#	238	0.00
35 RT	Benzene	10.000	10.441	-4.4	97	0.00
36 RT	1,2-Dichloroethane	10.000	10.619	-6.2	99	0.00
37 RT	Trichloroethene	10.000	10.429	-4.3	98	0.00
38 Rt	1,2-Dichloropropane	10.000	10.654	-6.5	99	0.00
39 t	Methacrylonitrile	10.000	10.468	-4.7	102	0.00
40 t	Methyl acrylate	10.000	10.949	-9.5	108	0.00
41 t	Tetrahydrofuran	20.000	52.881	-164.4#	290	0.00
42 t	1-Chlorobutane	10.000	10.349	-3.5	96	0.00
43 T	Dibromomethane	10.000	10.804	-8.0	100	0.00
44 T	Bromodichloromethane	10.000	10.973	-9.7	101	0.00
45 T	4-Methyl-2-Pentanone	50.000	52.867	-5.7	97	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	13.751	31.2#	77	0.00
47 t	Methyl methacrylate	20.000	10.476	47.6#	50	0.00
48 t	Ethyl methacrylate	10.000	10.759	-7.6	97	0.00
49 Rt	Toluene	10.000	10.505	-5.1	96	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
50 T	t-1,3-Dichloropropene	10.000	11.065	-10.6	100	0.00
51 T	cis-1,3-Dichloropropene	10.000	11.158	-11.6	102	0.00
52 RT	1,1,2-Trichloroethane	10.000	10.811	-8.1	101	0.00
53 t	1,3-Dichloropropane	10.000	10.558	-5.6	97	0.00
54 t	2-Hexanone	50.000	57.612	-15.2	108	0.00
55 t	Dibromochloromethane	10.000	10.920	-9.2	98	0.00
56 T	1,2-Dibromoethane	10.000	10.644	-6.4	97	0.00
57 S	4-Bromofluorobenzene	1.000	1.023	-2.3	100	0.00
58 RT	Tetrachloroethene	10.000	10.552	-5.5	98	0.00
59 Rt	Chlorobenzene	10.000	10.397	-4.0	96	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.992	-9.9	100	0.00
61 t	Pentachloroethane	10.000	11.630	-16.3	102	0.00
62 t	Hexachloroethane	10.000	11.152	-11.5	98	0.00
63 Rt	Ethyl Benzene	10.000	10.682	-6.8	98	0.00
64 RT	m/p-Xylenes	20.000	21.969	-9.8	100	0.00
65 RT	o-Xylene	10.000	11.033	-10.3	100	0.00
66 RT	Styrene	10.000	11.081	-10.8	100	0.00
67 t	Bromoform	10.000	11.318	-13.2	101	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.013	-1.3	98	0.00
69 T	Isopropylbenzene	10.000	10.913	-9.1	100	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	10.655	-6.5	99	0.00
71 T	1,2,3-Trichloropropane	10.000	9.396	6.0	82	0.00
72 t	Bromobenzene	10.000	10.840	-8.4	99	0.00
73 t	n-propylbenzene	10.000	11.200	-12.0	101	0.00
74 t	2-Chlorotoluene	10.000	11.043	-10.4	100	0.00
75 t	1,3,5-Trimethylbenzene	10.000	11.211	-12.1	102	0.00
76 t	4-Chlorotoluene	10.000	11.030	-10.3	100	0.00
77 t	tert-Butylbenzene	10.000	11.094	-10.9	101	0.00
78 t	1,2,4-Trimethylbenzene	10.000	10.964	-9.6	101	0.00
79 t	sec-Butylbenzene	10.000	11.245	-12.4	101	0.00
80	Nitrobenzene	50.000	10.515	79.0#	19	0.00
81 t	p-Isopropyltoluene	10.000	11.178	-11.8	101	0.00
82 t	1,3-Dichlorobenzene	10.000	11.210	-12.1	103	0.00
83 Rt	1,4-Dichlorobenzene	10.000	11.148	-11.5	103	0.00
84 t	n-Butylbenzene	10.000	11.250	-12.5	101	0.00
85 Rt	1,2-Dichlorobenzene	10.000	11.061	-10.6	102	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	10.339	-3.4	96	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	11.101	-11.0	102	0.00
88 t	Hexachlorobutadiene	10.000	11.480	-14.8	105	0.00
89 t	Naphthalene	10.000	10.788	-7.9	100	0.00
90 t	1,2,3-Trichlorobenzene	10.000	11.251	-12.5	101	0.00

(#) = Out of Range

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