

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU072419\
 Data File : VU033386.D
 Acq On : 24 Jul 2019 11:01
 Operator : JC/SP
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Jul 25 07:52:24 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA U\METHOD\524U072219DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Wed Jul 24 03:53:19 2019
 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	10.243	-2.4	101	0.00
3 t	Chloromethane	10.000	8.875	11.3	90	0.00
4 Rt	Vinyl Chloride	10.000	9.272	7.3	92	0.00
5 T	Bromomethane	10.000	9.240	7.6	96	0.00
6 T	Chloroethane	10.000	8.423	15.8	78	0.00
7 T	Trichlorofluoromethane	10.000	9.770	2.3	98	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.807	1.9	100	0.00
9 Rt	1,1-Dichloroethene	10.000	9.552	4.5	96	0.00
10 t	Iodomethane	10.000	7.742	22.6	73	0.00
11 t	Allyl Chloride	10.000	9.072	9.3	91	0.00
12 t	Acrylonitrile	20.000	15.859	20.7	83	0.00
13 T	Acetone	51.000	49.373	3.2	106	-0.02
14 T	Carbon Disulfide	10.000	9.099	9.0	93	0.00
15 RT	Methylene Chloride	10.000	8.431	15.7	94	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.024	9.8	93	0.00
17 t	1,1-Dichloroethane	10.000	12.616	-26.2	129	0.00
18 T	2-Butanone	50.000	52.395	-4.8	96	-0.01
19	Cyclohexane	10.000	10.799	-8.0	98	0.00
20	Methylcyclohexane	10.000	11.399	-14.0	100	0.00
21 T	2,2-Dichloropropane	10.000	10.496	-5.0	112	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.007	-0.1	97	0.00
23 t	Diethyl Ether	10.000	8.942	10.6	87	0.00
24 t	tert-Butyl Alcohol	100.000	77.509	22.5	72	-0.05
25 t	Methyl tert-Butyl Ether	10.000	8.740	12.6	86	0.00
26 t	Bromochloromethane	10.000	9.839	1.6	97	0.00
27 t	Chloroform	10.000	9.135	8.7	100	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.145	-1.4	100	0.00
29 T	1,1-Dichloropropene	10.000	10.515	-5.2	99	0.00
30 RT	Carbon Tetrachloride	10.000	10.175	-1.8	100	0.00
31 t	Isopropyl Ether	10.000	10.165	-1.6	98	0.00
32	Ethyl-t-butyl ether	10.000	10.207	-2.1	96	0.00
33	Tert-Amyl methyl ether	10.000	10.400	-4.0	91	0.00
34 t	Propionitrile	50.000	47.963	4.1	83	-0.02
35 RT	Benzene	10.000	10.429	-4.3	99	0.00
36 RT	1,2-Dichloroethane	10.000	9.875	1.3	96	0.00
37 RT	Trichloroethene	10.000	10.209	-2.1	98	0.00
38 Rt	1,2-Dichloropropane	10.000	10.248	-2.5	99	0.00
39 t	Methacrylonitrile	10.000	9.392	6.1	86	0.00
40 t	Methyl acrylate	10.000	9.895	1.1	86	0.00
41 t	Tetrahydrofuran	20.000	19.146	4.3	83	-0.01
42 t	1-Chlorobutane	10.000	10.481	-4.8	97	0.00
43 T	Dibromomethane	10.000	9.514	4.9	92	0.00
44 T	Bromodichloromethane	10.000	10.015	-0.2	99	0.00
45 T	4-Methyl-2-Pentanone	50.000	50.324	-0.6	87	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	19.210	3.9	80	0.00

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Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
47 t	Methyl methacrylate	20.000	20.963	-4.8	89	0.00
48 t	Ethyl methacrylate	10.000	10.728	-7.3	91	0.00
49 Rt	Toluene	10.000	11.601	-16.0	102	0.00
50 T	t-1,3-Dichloropropene	10.000	10.703	-7.0	97	0.00
51 T	cis-1,3-Dichloropropene	10.000	10.365	-3.7	97	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.838	1.6	94	0.00
53 t	1,3-Dichloropropane	10.000	10.129	-1.3	96	0.00
54 t	2-Hexanone	50.000	53.315	-6.6	92	0.00
55 t	Dibromochloromethane	10.000	10.320	-3.2	97	0.00
56 T	1,2-Dibromoethane	10.000	9.773	2.3	92	0.00
57 S	4-Bromofluorobenzene	1.000	1.279	-27.9	113	0.00
58 RT	Tetrachloroethene	10.000	11.044	-10.4	105	0.00
59 Rt	Chlorobenzene	10.000	11.085	-10.9	102	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.500	-5.0	103	0.00
61 t	Pentachloroethane	10.000	10.698	-7.0	106	0.00
62 t	Hexachloroethane	10.000	10.895	-8.9	104	0.00
63 Rt	Ethyl Benzene	10.000	11.908	-19.1	101	0.00
64 RT	m/p-Xylenes	20.000	24.834	-24.2	106	0.00
65 RT	o-Xylene	10.000	11.914	-19.1	102	0.00
66 RT	Styrene	10.000	12.449	-24.5	102	0.00
67 t	Bromoform	10.000	10.618	-6.2	97	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.044	-4.4	98	0.00
69 T	Isopropylbenzene	10.000	12.110	-21.1	103	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.806	1.9	95	0.00
71 T	1,2,3-Trichloropropane	10.000	9.738	2.6	98	0.00
72 t	Bromobenzene	10.000	11.175	-11.8	101	0.00
73 t	n-propylbenzene	10.000	12.652	-26.5	106	0.00
74 t	2-Chlorotoluene	10.000	12.244	-22.4	106	0.00
75 t	1,3,5-Trimethylbenzene	10.000	12.779	-27.8	105	0.00
76 t	4-Chlorotoluene	10.000	12.398	-24.0	107	0.00
77 t	tert-Butylbenzene	10.000	12.033	-20.3	105	0.00
78 t	1,2,4-Trimethylbenzene	10.000	12.904	-29.0	106	0.00
79 t	sec-Butylbenzene	10.000	12.390	-23.9	105	0.00
80	Nitrobenzene	50.000	32.474	35.1#	80	0.00
81 t	p-Isopropyltoluene	10.000	12.839	-28.4	106	0.00
82 t	1,3-Dichlorobenzene	10.000	11.339	-13.4	105	0.00
83 Rt	1,4-Dichlorobenzene	10.000	11.596	-16.0	105	0.00
84 t	n-Butylbenzene	10.000	12.722	-27.2	108	0.00
85 Rt	1,2-Dichlorobenzene	10.000	11.152	-11.5	104	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	10.045	-0.4	93	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	11.697	-17.0	100	0.00
88 t	Hexachlorobutadiene	10.000	11.323	-13.2	110	0.00
89 t	Naphthalene	10.000	8.915	10.9	87	0.00
90 t	1,2,3-Trichlorobenzene	10.000	11.413	-14.1	98	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0				