

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101823\
 Data File : VU055806.D
 Acq On : 18 Oct 2023 18:19
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Oct 19 03:42:24 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U101823DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Oct 19 03:10:04 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	94	0.00
2 T	Dichlorodifluoromethane	10.000	9.495	5.1	81	0.00
3 t	Chloromethane	10.000	9.991	0.1	87	0.00
4 Rt	Vinyl Chloride	10.000	10.000	0.0	87	0.00
5 T	Bromomethane	10.000	10.097	-1.0	86	0.00
6 T	Chloroethane	10.000	10.202	-2.0	85	0.00
7 T	Trichlorofluoromethane	10.000	9.865	1.3	85	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.527	4.7	81	0.00
9 Rt	1,1-Dichloroethene	10.000	9.592	4.1	82	0.00
10 t	Iodomethane	10.000	10.190	-1.9	82	0.00
11 t	Allyl Chloride	10.000	9.763	2.4	81	0.00
12 t	Acrylonitrile	20.000	20.908	-4.5	85	0.00
13 T	Acetone	50.000	41.000	18.0	71	0.00
14 T	Carbon Disulfide	10.000	9.521	4.8	81	0.00
15 RT	Methylene Chloride	10.000	9.777	2.2	88	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.906	0.9	83	0.00
17 t	1,1-Dichloroethane	10.000	9.742	2.6	83	0.00
18 T	2-Butanone	50.000	43.044	13.9	70	0.00
19	Cyclohexane	10.000	10.227	-2.3	81	0.00
20	Methylcyclohexane	10.000	10.354	-3.5	79	0.00
21 T	2,2-Dichloropropane	10.000	9.277	7.2	79	0.00
22 RT	cis-1,2-Dichloroethene	10.000	9.799	2.0	80	0.00
23 t	Diethyl Ether	10.000	10.373	-3.7	89	0.00
24 t	tert-Butyl Alcohol	100.000	106.442	-6.4	93	0.03
25 t	Methyl tert-Butyl Ether	10.000	10.373	-3.7	84	0.00
26 t	Bromochloromethane	10.000	9.718	2.8	83	0.00
27 t	Chloroform	10.000	9.650	3.5	83	0.00
28 RT	1,1,1-Trichloroethane	10.000	9.790	2.1	83	0.00
29 T	1,1-Dichloropropene	10.000	9.993	0.1	82	0.00
30 RT	Carbon Tetrachloride	10.000	9.793	2.1	83	0.00
31 t	Isopropyl Ether	10.000	10.362	-3.6	83	0.00
32	Ethyl-t-butyl ether	10.000	10.364	-3.6	83	0.00
33	Tert-Amyl methyl ether	10.000	10.165	-1.6	81	0.00
34 t	Propionitrile	50.000	55.722	-11.4	90	0.00
35 RT	Benzene	10.000	9.615	3.8	81	0.00
36 RT	1,2-Dichloroethane	10.000	9.612	3.9	83	0.00
37 RT	Trichloroethene	10.000	9.789	2.1	83	0.00
38 Rt	1,2-Dichloropropane	10.000	10.015	-0.2	85	0.00
39 t	Methacrylonitrile	10.000	10.738	-7.4	85	0.00
40 t	Methyl acrylate	10.000	8.855	11.4	78	0.00
41 t	Tetrahydrofuran	20.000	19.154	4.2	84	0.00
42 t	1-Chlorobutane	10.000	10.204	-2.0	83	0.00
43 T	Dibromomethane	10.000	9.844	1.6	86	0.00
44 T	Bromodichloromethane	10.000	9.778	2.2	83	0.00
45 T	4-Methyl-2-Pentanone	50.000	50.311	-0.6	81	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	22.794	-14.0	104	0.00
47 t	Methyl methacrylate	20.000	21.842	-9.2	84	0.00
48 t	Ethyl methacrylate	10.000	8.935	10.6	83	0.00
49 Rt	Toluene	10.000	10.338	-3.4	83	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
50 T	t-1,3-Dichloropropene	10.000	10.155	-1.5	84	0.00
51 T	cis-1,3-Dichloropropene	10.000	10.043	-0.4	82	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.716	2.8	86	0.00
53 t	1,3-Dichloropropane	10.000	9.909	0.9	86	0.00
54 t	2-Hexanone	50.000	43.644	12.7	69	0.00
55 t	Dibromochloromethane	10.000	9.709	2.9	84	0.00
56 T	1,2-Dibromoethane	10.000	10.028	-0.3	86	0.00
57 S	4-Bromofluorobenzene	1.000	0.961	3.9	91	0.00
58 RT	Tetrachloroethene	10.000	9.376	6.2	84	0.00
59 Rt	Chlorobenzene	10.000	10.019	-0.2	85	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	9.908	0.9	85	0.00
61 t	Pentachloroethane	10.000	9.310	6.9	78	0.00
62 t	Hexachloroethane	10.000	8.822	11.8	77	0.00
63 Rt	Ethyl Benzene	10.000	10.551	-5.5	83	0.00
64 RT	m/p-Xylenes	20.000	21.087	-5.4	83	0.00
65 RT	o-Xylene	10.000	9.837	1.6	79	0.00
66 RT	Styrene	10.000	10.305	-3.0	79	0.00
67 t	Bromoform	10.000	9.678	3.2	81	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.001	-0.1	90	0.00
69 T	Isopropylbenzene	10.000	10.128	-1.3	79	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.711	2.9	83	0.00
71 T	1,2,3-Trichloropropane	10.000	7.826	21.7	68	0.00
72 t	Bromobenzene	10.000	9.505	4.9	80	0.00
73 t	n-propylbenzene	10.000	8.879	11.2	80	0.00
74 t	2-Chlorotoluene	10.000	9.951	0.5	81	0.00
75 t	1,3,5-Trimethylbenzene	10.000	8.839	11.6	80	0.00
76 t	4-Chlorotoluene	10.000	9.979	0.2	79	0.00
77 t	tert-Butylbenzene	10.000	8.731	12.7	79	0.00
78 t	1,2,4-Trimethylbenzene	10.000	8.808	11.9	78	0.00
79 t	sec-Butylbenzene	10.000	8.740	12.6	78	0.00
80	Nitrobenzene	50.000	44.612	10.8	81	0.00
81 t	p-Isopropyltoluene	10.000	8.700	13.0	76	0.00
82 t	1,3-Dichlorobenzene	10.000	9.791	2.1	80	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.051	-0.5	81	0.00
84 t	n-Butylbenzene	10.000	8.535	14.6	74	0.00
85 Rt	1,2-Dichlorobenzene	10.000	9.913	0.9	81	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	11.218	-12.2	83	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	8.859	11.4	77	0.00
88 t	Hexachlorobutadiene	10.000	9.795	2.1	74	0.00
89 t	Naphthalene	10.000	8.412	15.9	75	0.00
90 t	1,2,3-Trichlorobenzene	10.000	8.867	11.3	78	0.00

(#) = Out of Range

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