

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101923\
 Data File : VU055808.D
 Acq On : 19 Oct 2023 09:30
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Oct 20 01:36:36 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	83	0.00
2 T	Dichlorodifluoromethane	10.000	9.194	8.1	69	0.00
3 t	Chloromethane	10.000	7.816	21.8	60	0.00
4 Rt	Vinyl Chloride	10.000	7.788	22.1	59	0.00
5 T	Bromomethane	10.000	8.812	11.9	66	0.00
6 T	Chloroethane	10.000	8.122	18.8	60	0.00
7 T	Trichlorofluoromethane	10.000	8.957	10.4	68	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.405	6.0	71	0.00
9 Rt	1,1-Dichloroethene	10.000	9.320	6.8	70	0.00
10 t	Iodomethane	10.000	10.039	-0.4	71	0.00
11 t	Allyl Chloride	10.000	9.238	7.6	68	0.00
12 t	Acrylonitrile	20.000	19.326	3.4	70	0.00
13 T	Acetone	50.000	39.201	21.6	60	0.00
14 T	Carbon Disulfide	10.000	9.359	6.4	70	0.00
15 RT	Methylene Chloride	10.000	9.021	9.8	72	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.706	2.9	71	0.00
17 t	1,1-Dichloroethane	10.000	9.437	5.6	71	0.00
18 T	2-Butanone	50.000	41.799	16.4	60	0.00
19	Cyclohexane	10.000	9.959	0.4	69	0.00
20	Methylcyclohexane	10.000	10.276	-2.8	69	0.00
21 T	2,2-Dichloropropane	10.000	9.540	4.6	72	0.00
22 RT	cis-1,2-Dichloroethene	10.000	9.726	2.7	70	0.00
23 t	Diethyl Ether	10.000	7.884	21.2	60	0.00
24 t	tert-Butyl Alcohol	100.000	90.176	9.8	69	0.01
25 t	Methyl tert-Butyl Ether	10.000	9.907	0.9	71	0.00
26 t	Bromochloromethane	10.000	9.655	3.5	72	0.00
27 t	Chloroform	10.000	9.538	4.6	72	0.00
28 RT	1,1,1-Trichloroethane	10.000	9.759	2.4	73	0.00
29 T	1,1-Dichloropropene	10.000	9.933	0.7	71	0.00
30 RT	Carbon Tetrachloride	10.000	9.769	2.3	73	0.00
31 t	Isopropyl Ether	10.000	9.717	2.8	69	0.00
32	Ethyl-t-butyl ether	10.000	9.926	0.7	70	0.00
33	Tert-Amyl methyl ether	10.000	9.970	0.3	70	0.00
34 t	Propionitrile	50.000	50.027	-0.1	71	0.00
35 RT	Benzene	10.000	9.639	3.6	71	0.00
36 RT	1,2-Dichloroethane	10.000	9.442	5.6	71	0.00
37 RT	Trichloroethene	10.000	9.499	5.0	71	0.00
38 Rt	1,2-Dichloropropane	10.000	9.433	5.7	71	0.00
39 t	Methacrylonitrile	10.000	9.975	0.3	69	0.00
40 t	Methyl acrylate	10.000	7.462	25.4	57	0.00
41 t	Tetrahydrofuran	20.000	18.188	9.1	70	0.00
42 t	1-Chlorobutane	10.000	9.862	1.4	71	0.00
43 T	Dibromomethane	10.000	9.356	6.4	72	0.00
44 T	Bromodichloromethane	10.000	9.613	3.9	72	0.00
45 T	4-Methyl-2-Pentanone	50.000	45.533	8.9	65	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	22.962	-14.8	92	0.00
47 t	Methyl methacrylate	20.000	20.579	-2.9	70	0.00
48 t	Ethyl methacrylate	10.000	8.471	15.3	69	0.00
49 Rt	Toluene	10.000	10.087	-0.9	71	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU101923\
 Data File : VU055808.D
 Acq On : 19 Oct 2023 09:30
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Oct 20 01:36:36 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)
50 T	t-1,3-Dichloropropene	10.000	9.826	1.7	71	0.00
51 T	cis-1,3-Dichloropropene	10.000	9.756	2.4	70	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.310	6.9	73	0.00
53 t	1,3-Dichloropropane	10.000	9.451	5.5	72	0.00
54 t	2-Hexanone	50.000	40.563	18.9	57	0.00
55 t	Dibromochloromethane	10.000	9.616	3.8	73	0.00
56 T	1,2-Dibromoethane	10.000	9.501	5.0	71	0.00
57 S	4-Bromofluorobenzene	1.000	1.239	-23.9	103	0.00
58 RT	Tetrachloroethene	10.000	9.296	7.0	74	0.00
59 Rt	Chlorobenzene	10.000	9.693	3.1	72	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	9.593	4.1	73	0.00
61 t	Pentachloroethane	10.000	9.947	0.5	73	0.00
62 t	Hexachloroethane	10.000	9.138	8.6	71	0.00
63 Rt	Ethyl Benzene	10.000	10.300	-3.0	71	0.00
64 RT	m/p-Xylenes	20.000	20.836	-4.2	72	0.00
65 RT	o-Xylene	10.000	10.162	-1.6	72	0.00
66 RT	Styrene	10.000	10.540	-5.4	71	0.00
67 t	Bromoform	10.000	9.853	1.5	73	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.018	-1.8	81	0.00
69 T	Isopropylbenzene	10.000	10.391	-3.9	71	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.379	6.2	71	0.00
71 T	1,2,3-Trichloropropane	10.000	8.699	13.0	66	0.00
72 t	Bromobenzene	10.000	9.666	3.3	71	0.00
73 t	n-propylbenzene	10.000	8.839	11.6	70	0.00
74 t	2-Chlorotoluene	10.000	10.070	-0.7	72	0.00
75 t	1,3,5-Trimethylbenzene	10.000	9.171	8.3	73	0.00
76 t	4-Chlorotoluene	10.000	10.291	-2.9	72	0.00
77 t	tert-Butylbenzene	10.000	8.859	11.4	71	0.00
78 t	1,2,4-Trimethylbenzene	10.000	8.965	10.4	70	0.00
79 t	sec-Butylbenzene	10.000	8.776	12.2	69	0.00
80	Nitrobenzene	50.000	38.979	22.0	62	0.00
81 t	p-Isopropyltoluene	10.000	8.727	12.7	67	0.00
82 t	1,3-Dichlorobenzene	10.000	9.879	1.2	71	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.131	-1.3	72	0.00
84 t	n-Butylbenzene	10.000	8.251	17.5	63	0.00
85 Rt	1,2-Dichlorobenzene	10.000	9.968	0.3	71	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	10.309	-3.1	67	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	8.670	13.3	66	0.00
88 t	Hexachlorobutadiene	10.000	10.620	-6.2	71	0.00
89 t	Naphthalene	10.000	8.324	16.8	66	0.00
90 t	1,2,3-Trichlorobenzene	10.000	9.030	9.7	70	0.00

(#) = Out of Range

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