

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU013023\
 Data File : VU052870.D
 Acq On : 30 Jan 2023 19:44
 Operator : JC/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Jan 31 13:21:57 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U012323DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jan 24 05:16:59 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	82	0.00
2 T	Dichlorodifluoromethane	10.000	8.448	15.5	67	0.00
3 t	Chloromethane	10.000	5.972	40.3#	51	0.00
4 Rt	Vinyl Chloride	10.000	6.922	30.8#	56	0.00
5 T	Bromomethane	10.000	9.602	4.0	75	0.00
6 T	Chloroethane	10.000	9.340	6.6	76	0.00
7 T	Trichlorofluoromethane	10.000	10.630	-6.3	86	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	11.249	-12.5	92	0.00
9 Rt	1,1-Dichloroethene	10.000	10.725	-7.2	86	0.00
10 t	Iodomethane	10.000	12.600	-26.0	91	0.00
11 t	Allyl Chloride	10.000	11.183	-11.8	89	0.00
12 t	Acrylonitrile	20.000	23.290	-16.4	90	0.00
13 T	Acetone	50.000	45.389	9.2	79	0.00
14 T	Carbon Disulfide	10.000	8.904	11.0	73	0.00
15 RT	Methylene Chloride	10.000	11.451	-14.5	103	0.00
16 RT	trans-1,2-Dichloroethene	10.000	10.654	-6.5	87	0.00
17 t	1,1-Dichloroethane	10.000	11.484	-14.8	94	0.00
18 T	2-Butanone	50.000	47.580	4.8	79	0.00
19	Cyclohexane	10.000	9.246	7.5	76	0.00
20	Methylcyclohexane	10.000	9.242	7.6	69	0.00
21 T	2,2-Dichloropropane	10.000	11.096	-11.0	92	0.00
22 RT	cis-1,2-Dichloroethene	10.000	11.080	-10.8	91	0.00
23 t	Diethyl Ether	10.000	11.125	-11.3	87	0.00
24 t	tert-Butyl Alcohol	100.000	127.278	-27.3	103	-0.02
25 t	Methyl tert-Butyl Ether	10.000	11.781	-17.8	94	0.00
26 t	Bromochloromethane	10.000	10.723	-7.2	90	0.00
27 t	Chloroform	10.000	12.034	-20.3	98	0.00
28 RT	1,1,1-Trichloroethane	10.000	12.010	-20.1	97	0.00
29 T	1,1-Dichloropropene	10.000	10.047	-0.5	81	0.00
30 RT	Carbon Tetrachloride	10.000	11.385	-13.8	90	0.00
31 t	Isopropyl Ether	10.000	11.096	-11.0	91	0.00
32	Ethyl-t-butyl ether	10.000	11.994	-19.9	95	0.00
33	Tert-Amyl methyl ether	10.000	10.932	-9.3	84	0.00
34 t	Propionitrile	50.000	62.631	-25.3	99	0.00
35 RT	Benzene	10.000	10.117	-1.2	81	0.00
36 RT	1,2-Dichloroethane	10.000	10.896	-9.0	88	0.00
37 RT	Trichloroethene	10.000	9.747	2.5	77	0.00
38 Rt	1,2-Dichloropropane	10.000	10.452	-4.5	82	0.00
39 t	Methacrylonitrile	10.000	10.990	-9.9	87	0.00
40 t	Methyl acrylate	10.000	10.554	-5.5	80	0.00
41 t	Tetrahydrofuran	20.000	21.649	-8.2	91	0.00
42 t	1-Chlorobutane	10.000	10.319	-3.2	84	0.00
43 T	Dibromomethane	10.000	10.548	-5.5	86	0.00
44 T	Bromodichloromethane	10.000	10.787	-7.9	88	0.00
45 T	4-Methyl-2-Pentanone	50.000	51.418	-2.8	76	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	18.211	8.9	73	0.00
47 t	Methyl methacrylate	20.000	21.266	-6.3	77	0.00
48 t	Ethyl methacrylate	10.000	9.934	0.7	70	0.00
49 Rt	Toluene	10.000	10.297	-3.0	79	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU013023\
 Data File : VU052870.D
 Acq On : 30 Jan 2023 19:44
 Operator : JC/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Jan 31 13:21:57 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U012323DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jan 24 05:16:59 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	9.854	1.5	74	0.00
51 T	cis-1,3-Dichloropropene	10.000	9.564	4.4	74	0.00
52 RT	1,1,2-Trichloroethane	10.000	10.780	-7.8	85	0.00
53 t	1,3-Dichloropropane	10.000	10.219	-2.2	80	0.00
54 t	2-Hexanone	50.000	44.236	11.5	66	0.00
55 t	Dibromochloromethane	10.000	10.440	-4.4	83	0.00
56 T	1,2-Dibromoethane	10.000	10.161	-1.6	80	0.00
57 S	4-Bromofluorobenzene	1.000	1.064	-6.4	82	0.00
58 RT	Tetrachloroethene	10.000	10.218	-2.2	81	0.00
59 Rt	Chlorobenzene	10.000	9.971	0.3	77	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.583	-5.8	85	0.00
61 t	Pentachloroethane	10.000	9.399	6.0	75	0.00
62 t	Hexachloroethane	10.000	9.135	8.7	70	0.00
63 Rt	Ethyl Benzene	10.000	10.415	-4.1	75	0.00
64 RT	m/p-Xylenes	20.000	22.616	-13.1	80	0.00
65 RT	o-Xylene	10.000	10.704	-7.0	77	0.00
66 RT	Styrene	10.000	9.659	3.4	81	0.00
67 t	Bromoform	10.000	10.360	-3.6	84	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.091	-9.1	83	0.00
69 T	Isopropylbenzene	10.000	10.704	-7.0	76	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	10.472	-4.7	82	0.00
71 T	1,2,3-Trichloropropane	10.000	5.902	41.0#	46	0.00
72 t	Bromobenzene	10.000	10.653	-6.5	82	0.00
73 t	n-propylbenzene	10.000	9.546	4.5	80	0.00
74 t	2-Chlorotoluene	10.000	11.374	-13.7	83	0.00
75 t	1,3,5-Trimethylbenzene	10.000	9.607	3.9	80	0.00
76 t	4-Chlorotoluene	10.000	11.304	-13.0	81	0.00
77 t	tert-Butylbenzene	10.000	10.053	-0.5	72	0.00
78 t	1,2,4-Trimethylbenzene	10.000	8.763	12.4	74	0.00
79 t	sec-Butylbenzene	10.000	8.406	15.9	70	0.00
80	Nitrobenzene	50.000	39.809	20.4	63	0.00
81 t	p-Isopropyltoluene	10.000	8.657	13.4	73	0.00
82 t	1,3-Dichlorobenzene	10.000	10.401	-4.0	79	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.397	-4.0	77	0.00
84 t	n-Butylbenzene	10.000	7.862	21.4	67	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.652	-6.5	80	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	8.828	11.7	70	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	10.104	-1.0	72	0.00
88 t	Hexachlorobutadiene	10.000	11.067	-10.7	85	0.00
89 t	Naphthalene	10.000	8.241	17.6	63	0.00
90 t	1,2,3-Trichlorobenzene	10.000	10.893	-8.9	78	0.00

(#) = Out of Range

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