

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032923\
 Data File : VU053476.D
 Acq On : 29 Mar 2023 20:33
 Operator : JC/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Mar 30 14:10:23 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U032923DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Mar 30 13:33:10 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	90	0.00
2 T	Dichlorodifluoromethane	10.000	9.856	1.4	83	0.00
3 t	Chloromethane	10.000	9.220	7.8	83	0.00
4 Rt	Vinyl Chloride	10.000	9.631	3.7	82	0.00
5 T	Bromomethane	10.000	10.323	-3.2	89	0.00
6 T	Chloroethane	10.000	9.419	5.8	86	0.00
7 T	Trichlorofluoromethane	10.000	9.160	8.4	81	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.322	6.8	84	0.00
9 Rt	1,1-Dichloroethene	10.000	9.597	4.0	86	0.00
10 t	Iodomethane	10.000	9.722	2.8	87	0.00
11 t	Allyl Chloride	10.000	10.007	-0.1	83	0.00
12 t	Acrylonitrile	20.000	19.361	3.2	83	0.00
13 T	Acetone	50.000	19.064	61.9#	35	-0.02
14 T	Carbon Disulfide	10.000	9.682	3.2	83	0.00
15 RT	Methylene Chloride	10.000	7.653	23.5	72	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.959	0.4	85	0.00
17 t	1,1-Dichloroethane	10.000	9.991	0.1	83	0.00
18 T	2-Butanone	50.000	29.987	40.0#	50	-0.03
19	Cyclohexane	10.000	9.820	1.8	82	0.00
20	Methylcyclohexane	10.000	10.005	-0.1	82	0.00
21 T	2,2-Dichloropropane	10.000	9.726	2.7	79	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.033	-0.3	83	0.00
23 t	Diethyl Ether	10.000	9.561	4.4	84	0.00
24 t	tert-Butyl Alcohol	100.000	97.315	2.7	81	-0.06
25 t	Methyl tert-Butyl Ether	10.000	10.227	-2.3	85	0.00
26 t	Bromochloromethane	10.000	9.957	0.4	83	0.00
27 t	Chloroform	10.000	10.049	-0.5	84	0.00
28 RT	1,1,1-Trichloroethane	10.000	9.966	0.3	81	0.00
29 T	1,1-Dichloropropene	10.000	10.251	-2.5	84	0.00
30 RT	Carbon Tetrachloride	10.000	10.454	-4.5	83	0.00
31 t	Isopropyl Ether	10.000	10.294	-2.9	85	-0.01
32	Ethyl-t-butyl ether	10.000	10.232	-2.3	84	0.00
33	Tert-Amyl methyl ether	10.000	10.413	-4.1	85	0.00
34 t	Propionitrile	50.000	50.114	-0.2	82	-0.03
35 RT	Benzene	10.000	10.011	-0.1	83	0.00
36 RT	1,2-Dichloroethane	10.000	9.940	0.6	85	0.00
37 RT	Trichloroethene	10.000	10.228	-2.3	82	0.00
38 Rt	1,2-Dichloropropane	10.000	10.232	-2.3	84	0.00
39 t	Methacrylonitrile	10.000	10.339	-3.4	87	0.00
40 t	Methyl acrylate	10.000	9.438	5.6	86	-0.01
41 t	Tetrahydrofuran	20.000	21.068	-5.3	85	-0.02
42 t	1-Chlorobutane	10.000	9.933	0.7	83	0.00
43 T	Dibromomethane	10.000	10.342	-3.4	84	0.00
44 T	Bromodichloromethane	10.000	10.314	-3.1	82	0.00
45 T	4-Methyl-2-Pentanone	50.000	49.807	0.4	81	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	18.696	6.5	82	0.00
47 t	Methyl methacrylate	20.000	21.087	-5.4	87	-0.01
48 t	Ethyl methacrylate	10.000	10.373	-3.7	82	0.00
49 Rt	Toluene	10.000	10.229	-2.3	83	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU032923\
 Data File : VU053476.D
 Acq On : 29 Mar 2023 20:33
 Operator : JC/MD
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Mar 30 14:10:23 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U032923DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Thu Mar 30 13:33:10 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	11.095	-11.0	85	0.00
51 T	cis-1,3-Dichloropropene	10.000	10.740	-7.4	84	0.00
52 RT	1,1,2-Trichloroethane	10.000	10.175	-1.8	84	0.00
53 t	1,3-Dichloropropane	10.000	10.193	-1.9	83	0.00
54 t	2-Hexanone	50.000	33.994	32.0#	56	0.00
55 t	Dibromochloromethane	10.000	10.802	-8.0	82	0.00
56 T	1,2-Dibromoethane	10.000	10.590	-5.9	85	0.00
57 S	4-Bromofluorobenzene	1.000	0.972	2.8	83	0.00
58 RT	Tetrachloroethene	10.000	10.407	-4.1	82	0.00
59 Rt	Chlorobenzene	10.000	10.147	-1.5	82	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.782	-7.8	83	0.00
61 t	Pentachloroethane	10.000	9.794	2.1	82	0.00
62 t	Hexachloroethane	10.000	9.365	6.3	81	0.00
63 Rt	Ethyl Benzene	10.000	10.540	-5.4	83	0.00
64 RT	m/p-Xylenes	20.000	21.215	-6.1	83	0.00
65 RT	o-Xylene	10.000	10.552	-5.5	83	0.00
66 RT	Styrene	10.000	10.954	-9.5	82	0.00
67 t	Bromoform	10.000	9.098	9.0	79	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.020	-2.0	87	0.00
69 T	Isopropylbenzene	10.000	10.545	-5.4	82	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	10.414	-4.1	85	0.00
71 T	1,2,3-Trichloropropane	10.000	11.366	-13.7	110	0.00
72 t	Bromobenzene	10.000	10.535	-5.4	84	0.00
73 t	n-propylbenzene	10.000	11.240	-12.4	85	0.00
74 t	2-Chlorotoluene	10.000	10.513	-5.1	83	0.00
75 t	1,3,5-Trimethylbenzene	10.000	10.908	-9.1	84	0.00
76 t	4-Chlorotoluene	10.000	10.630	-6.3	84	0.00
77 t	tert-Butylbenzene	10.000	10.595	-6.0	83	0.00
78 t	1,2,4-Trimethylbenzene	10.000	10.820	-8.2	83	0.00
79 t	sec-Butylbenzene	10.000	10.696	-7.0	82	0.00
80	Nitrobenzene	50.000	55.086	-10.2	78	0.00
81 t	p-Isopropyltoluene	10.000	10.938	-9.4	82	0.00
82 t	1,3-Dichlorobenzene	10.000	10.326	-3.3	82	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.403	-4.0	83	0.00
84 t	n-Butylbenzene	10.000	10.862	-8.6	80	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.539	-5.4	85	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	11.384	-13.8	90	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	10.817	-8.2	82	0.00
88 t	Hexachlorobutadiene	10.000	10.542	-5.4	81	0.00
89 t	Naphthalene	10.000	10.605	-6.1	80	0.00
90 t	1,2,3-Trichlorobenzene	10.000	11.174	-11.7	83	0.00

(#) = Out of Range

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