

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062325\
 Data File : VU063428.D
 Acq On : 23 Jun 2025 17:06
 Operator : MD/SY
 Sample : VSTDICV010
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICVVU062325

Quant Time: Jun 24 05:15:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U062325DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jun 24 05:06:50 2025
 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	103	0.00
2 T	Dichlorodifluoromethane	10.000	6.715	32.9#	68	0.00
3 t	Chloromethane	10.000	7.901	21.0	85	0.00
4 Rt	Vinyl Chloride	10.000	7.739	22.6	78	0.00
5 T	Bromomethane	10.000	8.107	18.9	87	0.00
6 T	Chloroethane	10.000	7.500	25.0	76	0.00
7 T	Trichlorofluoromethane	10.000	8.663	13.4	88	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.419	5.8	96	0.00
9 Rt	1,1-Dichloroethene	10.000	9.980	0.2	101	0.00
10 t	Iodomethane	10.000	12.484	-24.8	111	0.00
11 t	Allyl Chloride	10.000	9.992	0.1	101	0.00
12 t	Acrylonitrile	20.000	53.073	-165.4#	275	0.00
13 T	Acetone	50.000	36.130	27.7	75	0.00
14 T	Carbon Disulfide	10.000	9.898	1.0	101	0.00
15 RT	Methylene Chloride	10.000	8.959	10.4	94	0.00
16 RT	trans-1,2-Dichloroethene	10.000	10.251	-2.5	105	0.00
17 t	1,1-Dichloroethane	10.000	10.430	-4.3	106	0.00
18 T	2-Butanone	50.000	46.022	8.0	95	0.00
19	Cyclohexane	10.000	10.860	-8.6	109	0.00
20	Methylcyclohexane	10.000	11.686	-16.9	105	0.00
21 T	2,2-Dichloropropane	10.000	10.051	-0.5	104	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.853	-8.5	113	0.00
23 t	Diethyl Ether	10.000	8.222	17.8	82	0.00
24 t	tert-Butyl Alcohol	100.000	45.896	54.1#	45	0.00
25 t	Methyl tert-Butyl Ether	10.000	10.755	-7.6	107	0.00
26 t	Bromochloromethane	10.000	10.416	-4.2	106	0.00
27 t	Chloroform	10.000	10.619	-6.2	109	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.602	-6.0	104	0.00
29 T	1,1-Dichloropropene	10.000	10.004	-0.0	99	0.00
30 RT	Carbon Tetrachloride	10.000	10.815	-8.1	105	0.00
31 t	Isopropyl Ether	10.000	10.692	-6.9	107	0.00
32	Ethyl-t-butyl ether	10.000	0.000	100.0#	0	-4.50#
33	Tert-Amyl methyl ether	10.000	0.000	100.0#	0	-5.94#
34 t	Propionitrile	50.000	101.506	-103.0#	220	0.00
35 RT	Benzene	10.000	10.784	-7.8	105	0.00
36 RT	1,2-Dichloroethane	10.000	9.701	3.0	96	0.00
37 RT	Trichloroethene	10.000	9.923	0.8	98	0.00
38 Rt	1,2-Dichloropropane	10.000	10.383	-3.8	100	0.00
39 t	Methacrylonitrile	10.000	8.371	16.3	86	0.00
40 t	Methyl acrylate	10.000	10.805	-8.0	98	0.00
41 t	Tetrahydrofuran	20.000	42.681	-113.4#	230	0.00
42 t	1-Chlorobutane	10.000	10.403	-4.0	103	0.00
43 T	Dibromomethane	10.000	10.351	-3.5	103	0.00
44 T	Bromodichloromethane	10.000	10.803	-8.0	108	0.00
45 T	4-Methyl-2-Pentanone	50.000	55.253	-10.5	100	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	11.012	44.9#	59	0.00
47 t	Methyl methacrylate	20.000	9.804	51.0#	48	0.00
48 t	Ethyl methacrylate	10.000	10.000	0.0	103	0.00
49 Rt	Toluene	10.000	11.797	-18.0	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU062325\
 Data File : VU063428.D
 Acq On : 23 Jun 2025 17:06
 Operator : MD/SY
 Sample : VSTDICV010
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 ICSVU062325

Quant Time: Jun 24 05:15:13 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U062325DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Tue Jun 24 05:06:50 2025
 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	11.863	-18.6	106	0.00
51 T	cis-1,3-Dichloropropene	10.000	11.290	-12.9	103	0.00
52 RT	1,1,2-Trichloroethane	10.000	9.874	1.3	98	0.00
53 t	1,3-Dichloropropane	10.000	10.175	-1.8	99	0.00
54 t	2-Hexanone	50.000	47.244	5.5	98	0.00
55 t	Dibromochloromethane	10.000	11.013	-10.1	106	0.00
56 T	1,2-Dibromoethane	10.000	10.159	-1.6	102	0.00
57 S	4-Bromofluorobenzene	1.000	1.095	-9.5	98	0.00
58 RT	Tetrachloroethene	10.000	10.482	-4.8	101	0.00
59 Rt	Chlorobenzene	10.000	11.414	-14.1	108	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	9.976	0.2	98	0.00
61 t	Pentachloroethane	10.000	10.467	-4.7	107	0.00
62 t	Hexachloroethane	10.000	10.828	-8.3	106	0.00
63 Rt	Ethyl Benzene	10.000	12.408	-24.1	108	0.00
64 RT	m/p-Xylenes	20.000	21.238	-6.2	109	0.00
65 RT	o-Xylene	10.000	12.240	-22.4	108	0.00
66 RT	Styrene	10.000	10.412	-4.1	107	0.00
67 t	Bromoform	10.000	10.701	-7.0	103	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	0.972	2.8	93	0.00
69 T	Isopropylbenzene	10.000	13.991	-39.9#	122	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.680	3.2	97	0.00
71 T	1,2,3-Trichloropropane	10.000	8.495	15.1	95	0.00
72 t	Bromobenzene	10.000	11.207	-12.1	105	0.00
73 t	n-propylbenzene	10.000	10.430	-4.3	108	0.00
74 t	2-Chlorotoluene	10.000	11.949	-19.5	107	0.00
75 t	1,3,5-Trimethylbenzene	10.000	10.587	-5.9	108	0.00
76 t	4-Chlorotoluene	10.000	12.465	-24.6	113	0.00
77 t	tert-Butylbenzene	10.000	11.997	-20.0	106	0.00
78 t	1,2,4-Trimethylbenzene	10.000	10.383	-3.8	107	0.00
79 t	sec-Butylbenzene	10.000	12.220	-22.2	107	0.00
80	Nitrobenzene	50.000	10.734	78.5#	21	0.00
81 t	p-Isopropyltoluene	10.000	10.473	-4.7	108	0.00
82 t	1,3-Dichlorobenzene	10.000	11.397	-14.0	109	0.00
83 Rt	1,4-Dichlorobenzene	10.000	11.345	-13.5	109	0.00
84 t	n-Butylbenzene	10.000	12.351	-23.5	108	0.00
85 Rt	1,2-Dichlorobenzene	10.000	11.216	-12.2	106	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	10.426	-4.3	97	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	12.757	-27.6	118	0.00
88 t	Hexachlorobutadiene	10.000	10.514	-5.1	102	0.00
89 t	Naphthalene	10.000	10.410	-4.1	112	0.00
90 t	1,2,3-Trichlorobenzene	10.000	11.438	-14.4	105	0.00

(#) = Out of Range

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