

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110725\
 Data File : VU063678.D
 Acq On : 07 Nov 2025 08:55
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Nov 08 02:30:48 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U103125DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Mon Nov 03 00:51:51 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	93	0.00
2 T	Dichlorodifluoromethane	10.000	10.391	-3.9	92	0.00
3 t	Chloromethane	10.000	9.862	1.4	92	0.00
4 Rt	Vinyl Chloride	10.000	10.648	-6.5	93	0.00
5 T	Bromomethane	10.000	10.269	-2.7	99	0.01
6 T	Chloroethane	10.000	10.797	-8.0	96	0.00
7 T	Trichlorofluoromethane	10.000	11.000	-10.0	96	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	11.257	-12.6	99	0.00
9 Rt	1,1-Dichloroethene	10.000	10.909	-9.1	96	0.00
10 t	Iodomethane	10.000	11.049	-10.5	94	0.00
11 t	Allyl Chloride	10.000	10.502	-5.0	92	0.00
12 t	Acrylonitrile	20.000	19.892	0.5	83	0.00
13 T	Acetone	50.000	58.903	-17.8	109	0.00
14 T	Carbon Disulfide	10.000	10.221	-2.2	92	0.00
15 RT	Methylene Chloride	10.000	9.596	4.0	95	0.00
16 RT	trans-1,2-Dichloroethene	10.000	10.828	-8.3	94	0.00
17 t	1,1-Dichloroethane	10.000	10.810	-8.1	96	0.00
18 T	2-Butanone	50.000	56.816	-13.6	101	0.00
19	Cyclohexane	10.000	10.793	-7.9	93	0.00
20	Methylcyclohexane	10.000	10.081	-0.8	88	0.00
21 T	2,2-Dichloropropane	10.000	10.697	-7.0	95	0.00
22 RT	cis-1,2-Dichloroethene	10.000	10.936	-9.4	95	0.00
23 t	Diethyl Ether	10.000	10.297	-3.0	88	0.00
24 t	tert-Butyl Alcohol	100.000	68.727	31.3#	66	-0.01
25 t	Methyl tert-Butyl Ether	10.000	9.968	0.3	87	0.00
26 t	Bromochloromethane	10.000	10.455	-4.6	92	0.00
27 t	Chloroform	10.000	10.912	-9.1	97	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.811	-8.1	95	0.00
29 T	1,1-Dichloropropene	10.000	11.078	-10.8	96	0.00
30 RT	Carbon Tetrachloride	10.000	10.954	-9.5	93	0.00
31 t	Isopropyl Ether	10.000	10.728	-7.3	94	0.00
32	Ethyl-t-butyl ether	10.000	9.997	0.0	89	0.00
33	Tert-Amyl methyl ether	10.000	10.065	-0.6	88	0.00
34 t	Propionitrile	50.000	47.644	4.7	78	0.00
35 RT	Benzene	10.000	10.871	-8.7	96	0.00
36 RT	1,2-Dichloroethane	10.000	10.569	-5.7	91	0.00
37 RT	Trichloroethene	10.000	10.370	-3.7	97	0.00
38 Rt	1,2-Dichloropropane	10.000	10.123	-1.2	89	0.00
39 t	Methacrylonitrile	10.000	8.663	13.4	82	0.00
40 t	Methyl acrylate	10.000	8.782	12.2	82	0.00
41 t	Tetrahydrofuran	20.000	19.453	2.7	80	0.00
42 t	1-Chlorobutane	10.000	11.258	-12.6	94	0.00
43 T	Dibromomethane	10.000	9.720	2.8	83	0.00
44 T	Bromodichloromethane	10.000	9.985	0.2	86	0.00
45 T	4-Methyl-2-Pentanone	50.000	46.291	7.4	78	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	20.157	-0.8	84	0.00
47 t	Methyl methacrylate	20.000	18.773	6.1	77	0.00
48 t	Ethyl methacrylate	10.000	9.422	5.8	77	0.00
49 Rt	Toluene	10.000	10.472	-4.7	91	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU110725\
 Data File : VU063678.D
 Acq On : 07 Nov 2025 08:55
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Nov 08 02:30:48 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U103125DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Mon Nov 03 00:51:51 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	9.042	9.6	79	0.00
51 T	cis-1,3-Dichloropropene	10.000	9.849	1.5	85	0.00
52 RT	1,1,2-Trichloroethane	10.000	10.243	-2.4	87	0.00
53 t	1,3-Dichloropropane	10.000	9.803	2.0	86	0.00
54 t	2-Hexanone	50.000	55.827	-11.7	95	0.00
55 t	Dibromochloromethane	10.000	9.691	3.1	82	0.00
56 T	1,2-Dibromoethane	10.000	9.573	4.3	84	0.00
57 S	4-Bromofluorobenzene	1.000	1.206	-20.6	106	0.00
58 RT	Tetrachloroethene	10.000	11.187	-11.9	95	0.00
59 Rt	Chlorobenzene	10.000	10.557	-5.6	93	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	9.948	0.5	88	0.00
61 t	Pentachloroethane	10.000	9.267	7.3	82	0.00
62 t	Hexachloroethane	10.000	9.830	1.7	83	0.00
63 Rt	Ethyl Benzene	10.000	10.395	-3.9	92	0.00
64 RT	m/p-Xylenes	20.000	20.943	-4.7	91	0.00
65 RT	o-Xylene	10.000	10.450	-4.5	90	0.00
66 RT	Styrene	10.000	10.424	-4.2	89	0.00
67 t	Bromoform	10.000	8.900	11.0	75	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	0.975	2.5	81	0.00
69 T	Isopropylbenzene	10.000	10.450	-4.5	91	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	9.419	5.8	80	0.00
71 T	1,2,3-Trichloropropane	10.000	8.764	12.4	81	0.00
72 t	Bromobenzene	10.000	10.370	-3.7	90	0.00
73 t	n-propylbenzene	10.000	10.648	-6.5	91	0.00
74 t	2-Chlorotoluene	10.000	10.621	-6.2	93	0.00
75 t	1,3,5-Trimethylbenzene	10.000	10.400	-4.0	90	0.00
76 t	4-Chlorotoluene	10.000	10.744	-7.4	93	0.00
77 t	tert-Butylbenzene	10.000	10.072	-0.7	88	0.00
78 t	1,2,4-Trimethylbenzene	10.000	10.681	-6.8	90	0.00
79 t	sec-Butylbenzene	10.000	10.385	-3.8	89	0.00
80	Nitrobenzene	50.000	21.250	57.5#	38	0.00
81 t	p-Isopropyltoluene	10.000	10.425	-4.3	89	0.00
82 t	1,3-Dichlorobenzene	10.000	10.656	-6.6	91	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.503	-5.0	90	0.00
84 t	n-Butylbenzene	10.000	10.552	-5.5	87	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.390	-3.9	89	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	8.792	12.1	70	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	10.193	-1.9	83	0.00
88 t	Hexachlorobutadiene	10.000	11.050	-10.5	94	0.00
89 t	Naphthalene	10.000	8.272	17.3	69	0.00
90 t	1,2,3-Trichlorobenzene	10.000	9.777	2.2	79	0.00

(#) = Out of Range

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