

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU091525\
 Data File : VU063571.D
 Acq On : 15 Sep 2025 08:48
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Sep 16 04:54:06 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\524U091225DW.M
 Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
 QLast Update : Sat Sep 13 03:31:19 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	102	0.00
2 T	Dichlorodifluoromethane	10.000	9.788	2.1	99	0.00
3 t	Chloromethane	10.000	9.264	7.4	96	0.00
4 Rt	Vinyl Chloride	10.000	9.823	1.8	99	0.00
5 T	Bromomethane	10.000	8.676	13.2	96	0.00
6 T	Chloroethane	10.000	9.563	4.4	99	0.00
7 T	Trichlorofluoromethane	10.000	9.758	2.4	97	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	9.520	4.8	99	0.00
9 Rt	1,1-Dichloroethene	10.000	9.613	3.9	97	0.00
10 t	Iodomethane	10.000	9.657	3.4	94	0.00
11 t	Allyl Chloride	10.000	10.191	-1.9	101	0.00
12 t	Acrylonitrile	20.000	19.829	0.9	100	0.00
13 T	Acetone	50.000	57.206	-14.4	142	0.00
14 T	Carbon Disulfide	10.000	10.687	-6.9	106	0.00
15 RT	Methylene Chloride	10.000	8.919	10.8	98	0.00
16 RT	trans-1,2-Dichloroethene	10.000	9.982	0.2	102	0.00
17 t	1,1-Dichloroethane	10.000	9.845	1.5	101	0.00
18 T	2-Butanone	50.000	55.792	-11.6	122	0.00
19	Cyclohexane	10.000	8.931	10.7	90	0.00
20	Methylcyclohexane	10.000	9.955	0.4	98	0.00
21 T	2,2-Dichloropropane	10.000	10.301	-3.0	104	0.00
22 RT	cis-1,2-Dichloroethene	10.000	9.784	2.2	99	0.00
23 t	Diethyl Ether	10.000	9.648	3.5	99	0.00
24 t	tert-Butyl Alcohol	100.000	92.870	7.1	99	0.00
25 t	Methyl tert-Butyl Ether	10.000	9.783	2.2	99	0.00
26 t	Bromochloromethane	10.000	9.906	0.9	100	0.00
27 t	Chloroform	10.000	9.634	3.7	99	0.00
28 RT	1,1,1-Trichloroethane	10.000	10.147	-1.5	101	0.00
29 T	1,1-Dichloropropene	10.000	9.788	2.1	102	0.00
30 RT	Carbon Tetrachloride	10.000	10.776	-7.8	103	0.00
31 t	Isopropyl Ether	10.000	9.711	2.9	100	0.00
32	Ethyl-t-butyl ether	10.000	9.687	3.1	101	0.00
33	Tert-Amyl methyl ether	10.000	9.804	2.0	102	0.00
34 t	Propionitrile	50.000	48.204	3.6	100	0.00
35 RT	Benzene	10.000	9.649	3.5	98	0.00
36 RT	1,2-Dichloroethane	10.000	9.658	3.4	99	0.00
37 RT	Trichloroethene	10.000	9.736	2.6	100	0.00
38 Rt	1,2-Dichloropropane	10.000	9.918	0.8	93	0.00
39 t	Methacrylonitrile	10.000	9.174	8.3	102	0.00
40 t	Methyl acrylate	10.000	9.204	8.0	97	0.00
41 t	Tetrahydrofuran	20.000	18.152	9.2	100	0.00
42 t	1-Chlorobutane	10.000	10.048	-0.5	102	0.00
43 T	Dibromomethane	10.000	10.189	-1.9	102	0.00
44 T	Bromodichloromethane	10.000	10.746	-7.5	102	0.00
45 T	4-Methyl-2-Pentanone	50.000	51.669	-3.3	102	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	12.898	35.5#	61	0.00
47 t	Methyl methacrylate	20.000	20.023	-0.1	99	0.00
48 t	Ethyl methacrylate	10.000	10.520	-5.2	100	0.00
49 Rt	Toluene	10.000	10.251	-2.5	99	0.00

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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	11.355	-13.6	108	0.00
51 T	cis-1,3-Dichloropropene	10.000	11.017	-10.2	102	0.00
52 RT	1,1,2-Trichloroethane	10.000	10.238	-2.4	100	0.00
53 t	1,3-Dichloropropane	10.000	10.023	-0.2	100	0.00
54 t	2-Hexanone	50.000	57.972	-15.9	118	0.00
55 t	Dibromochloromethane	10.000	11.278	-12.8	106	0.00
56 T	1,2-Dibromoethane	10.000	9.980	0.2	99	0.00
57 S	4-Bromofluorobenzene	1.000	1.278	-27.8	126	0.00
58 RT	Tetrachloroethene	10.000	10.254	-2.5	103	0.00
59 Rt	Chlorobenzene	10.000	10.189	-1.9	99	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	10.769	-7.7	102	0.00
61 t	Pentachloroethane	10.000	10.375	-3.8	99	0.00
62 t	Hexachloroethane	10.000	11.596	-16.0	105	0.00
63 Rt	Ethyl Benzene	10.000	10.345	-3.5	98	0.00
64 RT	m/p-Xylenes	20.000	20.708	-3.5	98	0.00
65 RT	o-Xylene	10.000	10.304	-3.0	99	0.00
66 RT	Styrene	10.000	10.614	-6.1	98	0.00
67 t	Bromoform	10.000	11.830	-18.3	106	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	1.037	-3.7	97	0.00
69 T	Isopropylbenzene	10.000	10.376	-3.8	98	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	10.058	-0.6	98	0.00
71 T	1,2,3-Trichloropropane	10.000	10.106	-1.1	102	0.00
72 t	Bromobenzene	10.000	10.291	-2.9	100	0.00
73 t	n-propylbenzene	10.000	10.688	-6.9	97	0.00
74 t	2-Chlorotoluene	10.000	10.241	-2.4	97	0.00
75 t	1,3,5-Trimethylbenzene	10.000	10.514	-5.1	98	0.00
76 t	4-Chlorotoluene	10.000	10.015	-0.2	97	0.00
77 t	tert-Butylbenzene	10.000	10.544	-5.4	100	0.00
78 t	1,2,4-Trimethylbenzene	10.000	10.818	-8.2	99	0.00
79 t	sec-Butylbenzene	10.000	10.491	-4.9	98	0.00
80	Nitrobenzene	50.000	46.363	7.3	99	0.00
81 t	p-Isopropyltoluene	10.000	10.803	-8.0	98	0.00
82 t	1,3-Dichlorobenzene	10.000	10.279	-2.8	98	0.00
83 Rt	1,4-Dichlorobenzene	10.000	10.071	-0.7	99	0.00
84 t	n-Butylbenzene	10.000	10.607	-6.1	96	0.00
85 Rt	1,2-Dichlorobenzene	10.000	10.124	-1.2	98	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	11.750	-17.5	107	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	10.269	-2.7	96	0.00
88 t	Hexachlorobutadiene	10.000	10.595	-6.0	100	0.00
89 t	Naphthalene	10.000	9.529	4.7	90	0.00
90 t	1,2,3-Trichlorobenzene	10.000	10.151	-1.5	95	0.00

(#) = Out of Range

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