

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU091525\
 Data File : VU063584.D
 Acq On : 15 Sep 2025 17:45
 Operator : MD/SY
 Sample : VSTDCCC010
 Misc : 25mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 LabSampleId :
 VSTDCCC010

Quant Time: Sep 16 04:59:09 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	111	0.00
2 T	Dichlorodifluoromethane	10.000	8.020	19.8	88	0.00
3 t	Chloromethane	10.000	7.810	21.9	88	0.00
4 Rt	Vinyl Chloride	10.000	7.995	20.0	88	0.00
5 T	Bromomethane	10.000	7.413	25.9	89	0.00
6 T	Chloroethane	10.000	7.642	23.6	86	0.00
7 T	Trichlorofluoromethane	10.000	8.008	19.9	87	0.00
8	1,1,2-Trichloro-1,2,2-trifl	10.000	7.936	20.6	90	0.00
9 Rt	1,1-Dichloroethene	10.000	7.835	21.6	87	0.00
10 t	Iodomethane	10.000	8.323	16.8	88	0.00
11 t	Allyl Chloride	10.000	7.670	23.3	83	0.00
12 t	Acrylonitrile	20.000	16.066	19.7	88	0.00
13 T	Acetone	50.000	23.945	52.1#	65	0.00
14 T	Carbon Disulfide	10.000	7.163	28.4	78	0.00
15 RT	Methylene Chloride	10.000	7.968	20.3	95	0.00
16 RT	trans-1,2-Dichloroethene	10.000	8.064	19.4	89	0.00
17 t	1,1-Dichloroethane	10.000	7.980	20.2	89	0.00
18 T	2-Butanone	50.000	31.927	36.1#	76	0.00
19	Cyclohexane	10.000	7.651	23.5	84	0.00
20	Methylcyclohexane	10.000	8.019	19.8	86	0.00
21 T	2,2-Dichloropropane	10.000	7.303	27.0	80	0.00
22 RT	cis-1,2-Dichloroethene	10.000	7.908	20.9	87	0.00
23 t	Diethyl Ether	10.000	7.954	20.5	89	0.00
24 t	tert-Butyl Alcohol	100.000	74.045	26.0	86	0.00
25 t	Methyl tert-Butyl Ether	10.000	7.928	20.7	87	0.00
26 t	Bromochloromethane	10.000	8.443	15.6	92	0.00
27 t	Chloroform	10.000	7.952	20.5	89	0.00
28 RT	1,1,1-Trichloroethane	10.000	8.165	18.4	89	0.00
29 T	1,1-Dichloropropene	10.000	8.415	15.9	96	0.00
30 RT	Carbon Tetrachloride	10.000	8.686	13.1	90	0.00
31 t	Isopropyl Ether	10.000	7.810	21.9	88	0.00
32	Ethyl-t-butyl ether	10.000	7.768	22.3	88	0.00
33	Tert-Amyl methyl ether	10.000	9.598	4.0	109	0.00
34 t	Propionitrile	50.000	39.815	20.4	90	0.00
35 RT	Benzene	10.000	9.188	8.1	101	0.00
36 RT	1,2-Dichloroethane	10.000	9.641	3.6	108	0.00
37 RT	Trichloroethene	10.000	8.143	18.6	91	0.00
38 Rt	1,2-Dichloropropane	10.000	8.617	13.8	88	0.00
39 t	Methacrylonitrile	10.000	7.562	24.4	92	0.00
40 t	Methyl acrylate	10.000	7.213	27.9	83	0.00
41 t	Tetrahydrofuran	20.000	14.422	27.9	86	0.00
42 t	1-Chlorobutane	10.000	8.844	11.6	98	0.00
43 T	Dibromomethane	10.000	8.382	16.2	92	0.00
44 T	Bromodichloromethane	10.000	8.701	13.0	90	0.00
45 T	4-Methyl-2-Pentanone	50.000	47.524	5.0	102	0.00
46 t	t-1,4-Dichloro-2-butene	20.000	13.433	32.8#	69	0.00
47 t	Methyl methacrylate	20.000	17.835	10.8	96	0.00
48 t	Ethyl methacrylate	10.000	8.646	13.5	90	0.00
49 Rt	Toluene	10.000	8.734	12.7	92	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	7.891	21.1	82	0.00
51 T	cis-1,3-Dichloropropene	10.000	9.250	7.5	94	0.00
52 RT	1,1,2-Trichloroethane	10.000	8.414	15.9	90	0.00
53 t	1,3-Dichloropropane	10.000	8.356	16.4	91	0.00
54 t	2-Hexanone	50.000	38.128	23.7	85	0.00
55 t	Dibromochloromethane	10.000	8.197	18.0	84	0.00
56 T	1,2-Dibromoethane	10.000	8.388	16.1	91	0.00
57 S	4-Bromofluorobenzene	1.000	0.846	15.4	91	0.00
58 RT	Tetrachloroethene	10.000	8.091	19.1	88	0.00
59 Rt	Chlorobenzene	10.000	8.412	15.9	89	0.00
60 T	1,1,1,2-Tetrachloroethane	10.000	8.299	17.0	85	0.00
61 t	Pentachloroethane	10.000	8.515	14.8	88	0.00
62 t	Hexachloroethane	10.000	8.770	12.3	86	0.00
63 Rt	Ethyl Benzene	10.000	8.524	14.8	88	0.00
64 RT	m/p-Xylenes	20.000	17.367	13.2	90	0.00
65 RT	o-Xylene	10.000	8.512	14.9	89	0.00
66 RT	Styrene	10.000	9.020	9.8	91	0.00
67 t	Bromoform	10.000	8.371	16.3	82	0.00
68 S	1,2-Dichlorobenzene-d4	1.000	0.924	7.6	94	0.00
69 T	Isopropylbenzene	10.000	8.581	14.2	88	0.00
70 T	1,1,2,2-Tetrachloroethane	10.000	8.587	14.1	91	0.00
71 T	1,2,3-Trichloropropane	10.000	8.950	10.5	98	0.00
72 t	Bromobenzene	10.000	8.729	12.7	92	0.00
73 t	n-propylbenzene	10.000	8.834	11.7	88	0.00
74 t	2-Chlorotoluene	10.000	8.794	12.1	91	0.00
75 t	1,3,5-Trimethylbenzene	10.000	8.947	10.5	91	0.00
76 t	4-Chlorotoluene	10.000	8.551	14.5	90	0.00
77 t	tert-Butylbenzene	10.000	8.725	12.8	90	0.00
78 t	1,2,4-Trimethylbenzene	10.000	9.167	8.3	91	0.00
79 t	sec-Butylbenzene	10.000	8.869	11.3	90	0.00
80	Nitrobenzene	50.000	35.214	29.6	80	0.00
81 t	p-Isopropyltoluene	10.000	9.104	9.0	90	0.00
82 t	1,3-Dichlorobenzene	10.000	8.721	12.8	90	0.00
83 Rt	1,4-Dichlorobenzene	10.000	8.506	14.9	91	0.00
84 t	n-Butylbenzene	10.000	9.010	9.9	89	0.00
85 Rt	1,2-Dichlorobenzene	10.000	8.719	12.8	92	0.00
86 t	1,2-Dibromo-3-Chloropropane	10.000	9.119	8.8	90	0.00
87 Rt	1,2,4-Trichlorobenzene	10.000	8.518	14.8	86	0.00
88 t	Hexachlorobutadiene	10.000	8.829	11.7	91	0.00
89 t	Naphthalene	10.000	8.656	13.4	89	0.00
90 t	1,2,3-Trichlorobenzene	10.000	9.197	8.0	94	0.00

(#) = Out of Range

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