

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091225\
 Data File : VN087811.D
 Acq On : 12 Sep 2025 10:04
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Sep 12 22:49:07 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N090425W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Sep 05 03:29:53 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	142	0.00
2 M	Dichlorodifluoromethane	1.865	1.614	13.5	131	0.00
3 M	Chloromethane	1.957	1.569	19.8	124	0.00
4 M	Vinyl Chloride	2.213	1.815	18.0	129	0.00
5 M	Bromomethane	1.298	0.863	33.5#	111	0.00
6 M	Chloroethane	1.512	1.146	24.2	120	0.00
7 M	Trichlorofluoromethane	3.243	2.751	15.2	135	0.00
8 T	Diethyl Ether	1.370	1.068	22.0	130	0.01
9	1,1,2-Trichlorotrifluoroeth	1.757	1.483	15.6	132	0.00
10 M	1,1-Dichloroethene	1.811	1.473	18.7	133	0.00
11	Methyl Iodide	1.282	0.747	41.7#	113	0.00
12	Methyl Acetate	2.792	2.322	16.8	130	0.00
13 M	Acrolein	0.497	0.474	4.6	156#	0.00
14 M	Acrylonitrile	1.202	0.967	19.6	131	-0.01
15 M	Acetone	0.375	0.317	15.5	141	-0.01
16 M	Carbon Disulfide	4.972	4.263	14.3	139	0.00
17	Allyl chloride	2.958	2.436	17.6	136	0.00
18 M	Methylene Chloride	2.052	1.712	16.6	138	0.00
19 M	trans-1,2-Dichloroethene	1.847	1.549	16.1	133	0.00
20 T	Diisopropyl ether	6.802	5.379	20.9	129	0.00
21 M	1,1-Dichloroethane	3.725	3.057	17.9	132	0.00
22 M	cis-1,2-Dichloroethene	2.242	1.822	18.7	130	0.00
23 M	tert-Butyl Alcohol	0.425	0.335	21.2	124	0.00
24 M	Methyl tert-Butyl Ether	6.466	5.329	17.6	137	0.00
25 M	Chloroform	3.899	3.225	17.3	131	0.00
26	Cyclohexane	2.837	2.344	17.4	136	-0.01
27 s	1,2-Dichloroethane-d4	2.563	2.485	3.0	141	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	143	0.00
29	1,1-Dichloropropene	0.488	0.405	17.0	130	0.00
30 M	2-Butanone	0.336	0.282	16.1	133	0.00
31	2,2-Dichloropropane	0.588	0.534	9.2	155#	0.00
32 M	1,1,1-Trichloroethane	0.618	0.524	15.2	129	0.00
33 M	Carbon Tetrachloride	0.515	0.434	15.7	134	0.00
34 M	Benzene	1.576	1.323	16.1	132	0.00
35	Methacrylonitrile	0.351	0.311	11.4	130	0.00
36 M	1,2-Dichloroethane	0.595	0.502	15.6	130	0.00
37 M	Trichloroethene	0.352	0.313	11.1	139	0.00
38	Methylcyclohexane	0.537	0.464	13.6	140	0.00
39 M	1,2-Dichloropropane	0.404	0.337	16.6	129	0.00
40	Dibromomethane	0.300	0.263	12.3	137	0.00
41 M	Bromodichloromethane	0.614	0.533	13.2	134	0.00
42 M	Vinyl Acetate	1.122	0.893	20.4	128	0.00
43	Ethyl Acetate	0.686	0.580	15.5	129	0.00
44	Isopropyl Acetate	1.076	0.867	19.4	125	0.00
45 T	1,4-Dioxane	0.007	0.006	14.3	131	0.00
46	Methyl methacrylate	0.496	0.338	31.9#	109	0.00
47	n-amyl Acetate	0.727	0.474	34.8#	131	-0.13
48 M	t-1,3-Dichloropropene	0.638	0.527	17.4	136	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091225\
 Data File : VN087811.D
 Acq On : 12 Sep 2025 10:04
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Sep 12 22:49:07 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N090425W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Sep 05 03:29:53 2025
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.653	0.549	15.9	139	0.00
50 M	1,1,2-Trichloroethane	0.388	0.342	11.9	139	0.00
51	Ethyl methacrylate	0.604	0.462	23.5	129	0.00
52	1,3-Dichloropropane	0.674	0.572	15.1	132	0.00
53 M	Dibromochloromethane	0.435	0.378	13.1	147	0.00
54 M	1,2-Dibromoethane	0.404	0.337	16.6	133	0.00
55 M	2-Chloroethyl vinyl ether	0.346	0.323	6.6	145	0.00
56 M	Bromoform	0.284	0.237	16.5	140	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	145	0.00
58 M	4-Methyl-2-Pentanone	0.709	0.587	17.2	130	0.00
59 M	2-Hexanone	0.464	0.371	20.0	128	0.00
60 S	4-Bromofluorobenzene	0.482	0.491	-1.9	145	0.00
61 M	Tetrachloroethene	0.299	0.271	9.4	140	0.00
62 M	Toluene	1.815	1.521	16.2	132	0.00
63 S	Toluene-d8	1.395	1.380	1.1	141	0.00
64 M	Chlorobenzene	1.114	0.953	14.5	137	0.00
65	1,1,1,2-Tetrachloroethane	0.392	0.350	10.7	143	0.00
66 M	Ethyl Benzene	1.958	1.644	16.0	136	0.00
67 M	m/p-Xylenes	0.727	0.608	16.4	137	0.00
68 M	o-Xylene	0.705	0.596	15.5	135	0.00
69 M	Styrene	1.225	0.998	18.5	134	0.00
70	Isopropylbenzene	1.787	1.513	15.3	141	0.00
71 M	1,1,2,2-Tetrachloroethane	0.642	0.544	15.3	130	0.00
72	1,2,3-Trichloropropane	0.631	0.511	19.0	116	0.00
73	Bromobenzene	0.437	0.374	14.4	137	0.00
74	n-propylbenzene	2.254	1.923	14.7	138	0.00
75	2-Chlorotoluene	1.357	1.133	16.5	132	0.00
76	1,3,5-Trimethylbenzene	1.509	1.280	15.2	137	0.00
77	t-1,4-Dichloro-2-butene	0.213	0.153	28.2#	121	0.00
78	4-Chlorotoluene	1.368	1.135	17.0	130	0.00
79	tert-butylbenzene	1.263	1.094	13.4	140	0.00
80	1,2,4-Trimethylbenzene	1.547	1.318	14.8	136	0.00
81	sec-Butylbenzene	1.858	1.631	12.2	141	0.00
82	p-Isopropyltoluene	1.531	1.348	12.0	143	0.00
83 M	1,3-Dichlorobenzene	0.843	0.748	11.3	137	0.00
84 M	1,4-Dichlorobenzene	0.845	0.753	10.9	140	0.00
85	n-Butylbenzene	1.472	1.323	10.1	149	0.00
86 T	Hexachloroethane	0.311	0.270	13.2	136	0.00
87 M	1,2-Dichlorobenzene	0.795	0.722	9.2	142	0.00
88	1,2-Dibromo-3-Chloropropane	0.143	0.113	21.0	117	0.00
89	1,2,4-Trichlorobenzene	0.452	0.402	11.1	145	0.00
90	Hexachlorobutadiene	0.152	0.163	-7.2	162#	0.00
91 M	Naphthalene	1.664	1.347	19.1	135	0.00
92	1,2,3-Trichlorobenzene	0.452	0.402	11.1	145	0.00

(#) = Out of Range

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