

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090425\
 Data File : VN087731.D
 Acq On : 04 Sep 2025 18:28
 Operator : JC\MD
 Sample : VSTDICV020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN090425

Quant Time: Sep 05 03:32:36 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	107	0.00
2 M	Dichlorodifluoromethane	1.865	1.713	8.2	105	0.00
3 M	Chloromethane	1.957	1.766	9.8	106	0.00
4 M	Vinyl Chloride	2.213	1.936	12.5	104	0.00
5 M	Bromomethane	1.298	1.158	10.8	112	0.01
6 M	Chloroethane	1.512	1.292	14.6	102	0.00
7 M	Trichlorofluoromethane	3.243	2.785	14.1	103	0.00
8 T	Diethyl Ether	1.370	1.207	11.9	111	0.00
9	1,1,2-Trichlorotrifluoroeth	1.757	1.510	14.1	102	0.00
10 M	1,1-Dichloroethene	1.811	1.584	12.5	108	0.00
11	Methyl Iodide	1.282	0.968	24.5	111	0.01
12	Methyl Acetate	2.792	2.549	8.7	108	0.00
13 M	Acrolein	0.497	0.453	8.9	113	0.00
14 M	Acrylonitrile	1.202	1.062	11.6	108	0.00
15 M	Acetone	0.375	0.317	15.5	107	0.00
16 M	Carbon Disulfide	4.972	4.458	10.3	110	0.00
17	Allyl chloride	2.958	2.524	14.7	107	0.00
18 M	Methylene Chloride	2.052	1.826	11.0	111	0.00
19 M	trans-1,2-Dichloroethene	1.847	1.624	12.1	105	0.00
20 T	Diisopropyl ether	6.802	5.657	16.8	102	0.00
21 M	1,1-Dichloroethane	3.725	3.283	11.9	107	0.00
22 M	cis-1,2-Dichloroethene	2.242	1.939	13.5	105	0.00
23 M	tert-Butyl Alcohol	0.425	0.382	10.1	107	-0.01
24 M	Methyl tert-Butyl Ether	6.466	5.592	13.5	108	0.00
25 M	Chloroform	3.899	3.382	13.3	104	0.00
26	Cyclohexane	2.837	2.462	13.2	108	0.00
27 s	1,2-Dichloroethane-d4	2.563	2.583	-0.8	110	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	111	0.00
29	1,1-Dichloropropene	0.488	0.417	14.5	104	0.00
30 M	2-Butanone	0.336	0.295	12.2	109	0.00
31	2,2-Dichloropropane	0.588	0.500	15.0	112	0.00
32 M	1,1,1-Trichloroethane	0.618	0.545	11.8	105	0.00
33 M	Carbon Tetrachloride	0.515	0.443	14.0	106	0.00
34 M	Benzene	1.576	1.392	11.7	108	0.00
35	Methacrylonitrile	0.351	0.326	7.1	105	0.00
36 M	1,2-Dichloroethane	0.595	0.523	12.1	105	0.00
37 M	Trichloroethene	0.352	0.292	17.0	101	0.00
38	Methylcyclohexane	0.537	0.449	16.4	105	0.00
39 M	1,2-Dichloropropane	0.404	0.348	13.9	104	0.00
40	Dibromomethane	0.300	0.269	10.3	109	0.00
41 M	Bromodichloromethane	0.614	0.533	13.2	104	0.00
42 M	Vinyl Acetate	1.122	0.990	11.8	110	0.00
43	Ethyl Acetate	0.686	0.594	13.4	102	0.00
44	Isopropyl Acetate	1.076	0.938	12.8	105	0.00
45 T	1,4-Dioxane	0.007	0.006	14.3	100	0.00
46	Methyl methacrylate	0.496	0.432	12.9	109	0.00
47	n-amyl Acetate	0.727	0.551	24.2	119	-0.10
48 M	t-1,3-Dichloropropene	0.638	0.545	14.6	109	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.653	0.563	13.8	111	0.00
50 M	1,1,2-Trichloroethane	0.388	0.342	11.9	108	0.00
51	Ethyl methacrylate	0.604	0.523	13.4	114	0.00
52	1,3-Dichloropropane	0.674	0.596	11.6	107	0.00
53 M	Dibromochloromethane	0.435	0.377	13.3	114	0.00
54 M	1,2-Dibromoethane	0.404	0.348	13.9	107	0.00
55 M	2-Chloroethyl vinyl ether	0.346	0.298	13.9	104	0.00
56 M	Bromoform	0.284	0.249	12.3	115	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	112	0.00
58 M	4-Methyl-2-Pentanone	0.709	0.625	11.8	107	0.00
59 M	2-Hexanone	0.464	0.395	14.9	105	0.00
60 S	4-Bromofluorobenzene	0.482	0.470	2.5	107	0.00
61 M	Tetrachloroethene	0.299	0.262	12.4	105	0.00
62 M	Toluene	1.815	1.618	10.9	108	0.00
63 S	Toluene-d8	1.395	1.380	1.1	109	0.00
64 M	Chlorobenzene	1.114	0.968	13.1	107	0.00
65	1,1,1,2-Tetrachloroethane	0.392	0.335	14.5	106	0.00
66 M	Ethyl Benzene	1.958	1.699	13.2	109	0.00
67 M	m/p-Xylenes	0.727	0.628	13.6	109	0.00
68 M	o-Xylene	0.705	0.612	13.2	107	0.00
69 M	Styrene	1.225	1.029	16.0	107	0.00
70	Isopropylbenzene	1.787	1.521	14.9	109	0.00
71 M	1,1,2,2-Tetrachloroethane	0.642	0.569	11.4	105	0.00
72	1,2,3-Trichloropropane	0.631	0.536	15.1	93	0.00
73	Bromobenzene	0.437	0.375	14.2	106	0.00
74	n-propylbenzene	2.254	1.932	14.3	107	0.00
75	2-Chlorotoluene	1.357	1.165	14.1	105	0.00
76	1,3,5-Trimethylbenzene	1.509	1.312	13.1	108	0.00
77	t-1,4-Dichloro-2-butene	0.213	0.162	23.9	99	0.00
78	4-Chlorotoluene	1.368	1.250	8.6	111	0.00
79	tert-butylbenzene	1.263	1.102	12.7	108	0.00
80	1,2,4-Trimethylbenzene	1.547	1.327	14.2	105	0.00
81	sec-Butylbenzene	1.858	1.615	13.1	107	0.00
82	p-Isopropyltoluene	1.531	1.327	13.3	109	0.00
83 M	1,3-Dichlorobenzene	0.843	0.719	14.7	102	0.00
84 M	1,4-Dichlorobenzene	0.845	0.730	13.6	105	0.00
85	n-Butylbenzene	1.472	1.280	13.0	111	0.00
86 T	Hexachloroethane	0.311	0.280	10.0	109	0.00
87 M	1,2-Dichlorobenzene	0.795	0.716	9.9	109	0.00
88	1,2-Dibromo-3-Chloropropane	0.143	0.130	9.1	104	0.00
89	1,2,4-Trichlorobenzene	0.452	0.396	12.4	110	0.00
90	Hexachlorobutadiene	0.152	0.138	9.2	106	0.00
91 M	Naphthalene	1.664	1.414	15.0	109	0.00
92	1,2,3-Trichlorobenzene	0.452	0.396	12.4	110	0.00

(#) = Out of Range

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