

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062322\
 Data File : VN073071.D
 Acq On : 24 Jun 2022 02:23
 Operator : JC\MD
 Sample : VSTDICV020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICSVN062322

Quant Time: Jun 24 05:01:00 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N062322W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Jun 24 04:54:59 2022
 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	100	0.00
2 M	Dichlorodifluoromethane	2.367	2.558	-8.1	101	0.00
3 M	Chloromethane	2.249	2.409	-7.1	105	0.00
4 M	Vinyl Chloride	2.013	2.164	-7.5	102	0.00
5 M	Bromomethane	0.663	0.945	-42.5#	113	0.00
6 M	Chloroethane	1.242	1.317	-6.0	99	0.00
7 M	Trichlorofluoromethane	3.654	3.800	-4.0	104	0.00
8 T	Diethyl Ether	1.495	1.580	-5.7	103	0.00
9	1,1,2-Trichlorotrifluoroeth	2.019	2.167	-7.3	107	0.00
10 M	1,1-Dichloroethene	2.066	2.118	-2.5	102	0.00
11	Methyl Iodide	1.971	1.658	15.9	93	0.00
12	Methyl Acetate	4.302	4.177	2.9	97	0.00
13 M	Acrolein	0.434	0.496	-14.3	110	0.00
14 M	Acrylonitrile	1.639	1.652	-0.8	100	0.00
15 M	Acetone	0.452	0.449	0.7	95	0.00
16 M	Carbon Disulfide	5.447	5.560	-2.1	105	0.00
17	Allyl chloride	3.921	4.146	-5.7	109	0.00
18 M	Methylene Chloride	2.464	2.514	-2.0	103	0.00
19 M	trans-1,2-Dichloroethene	2.180	2.287	-4.9	104	0.00
20 T	Diisopropyl ether	8.250	8.499	-3.0	104	0.00
21 M	1,1-Dichloroethane	4.329	4.385	-1.3	102	0.00
22 M	cis-1,2-Dichloroethene	2.592	2.660	-2.6	104	0.00
23 M	tert-Butyl Alcohol	0.669	0.666	0.4	103	0.00
24 M	Methyl tert-Butyl Ether	7.706	7.896	-2.5	102	0.00
25 M	Chloroform	4.308	4.431	-2.9	103	0.00
26	Cyclohexane	3.790	4.010	-5.8	107	0.00
27 s	1,2-Dichloroethane-d4	2.892	2.957	-2.2	100	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	98	0.00
29	1,1-Dichloropropene	0.542	0.562	-3.7	105	0.00
30 M	2-Butanone	0.418	0.425	-1.7	99	0.00
31	2,2-Dichloropropane	0.265	0.455	-71.7#	164#	0.00
32 M	1,1,1-Trichloroethane	0.666	0.701	-5.3	106	0.00
33 M	Carbon Tetrachloride	0.564	0.587	-4.1	105	0.00
34 M	Benzene	1.688	1.767	-4.7	102	0.00
35	Methacrylonitrile	0.372	0.396	-6.5	103	0.00
36 M	1,2-Dichloroethane	0.626	0.636	-1.6	101	0.00
37 M	Trichloroethene	0.420	0.426	-1.4	103	0.00
38	Methylcyclohexane	0.642	0.670	-4.4	104	0.00
39 M	1,2-Dichloropropane	0.437	0.455	-4.1	103	0.00
40	Dibromomethane	0.295	0.301	-2.0	100	0.00
41 M	Bromodichloromethane	0.594	0.604	-1.7	104	0.00
42 M	Vinyl Acetate	1.292	1.359	-5.2	105	0.00
43	Ethyl Acetate	0.787	0.854	-8.5	100	0.00
44	Isopropyl Acetate	1.215	1.246	-2.6	104	0.00
45 T	1,4-Dioxane	0.011	0.011	0.0	105	0.00
46	Methyl methacrylate	0.553	0.551	0.4	103	0.00
47	n-amyl Acetate	0.974	0.981	-0.7	106	0.00
48 M	t-1,3-Dichloropropene	0.556	0.614	-10.4	116	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.602	0.671	-11.5	114	0.00
50 M	1,1,2-Trichloroethane	0.428	0.442	-3.3	102	0.00
51	Ethyl methacrylate	0.725	0.733	-1.1	102	0.00
52	1,3-Dichloropropane	0.736	0.760	-3.3	104	0.00
53 M	Dibromochloromethane	0.445	0.441	0.9	104	0.00
54 M	1,2-Dibromoethane	0.449	0.462	-2.9	104	0.00
55 M	2-Chloroethyl vinyl ether	0.310	0.319	-2.9	100	0.00
56 M	Bromoform	0.326	0.314	3.7	105	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	98	0.00
58 M	4-Methyl-2-Pentanone	0.873	0.920	-5.4	102	0.00
59 M	2-Hexanone	0.666	0.701	-5.3	102	0.00
60 S	4-Bromofluorobenzene	0.574	0.573	0.2	102	0.00
61 M	Tetrachloroethene	0.388	0.407	-4.9	105	0.00
62 M	Toluene	1.994	2.102	-5.4	103	0.00
63 S	Toluene-d8	1.625	1.675	-3.1	100	0.00
64 M	Chlorobenzene	1.200	1.238	-3.2	103	0.00
65	1,1,1,2-Tetrachloroethane	0.456	0.466	-2.2	102	0.00
66 M	Ethyl Benzene	2.197	2.282	-3.9	104	0.00
67 M	m/p-Xylenes	0.824	0.872	-5.8	105	0.00
68 M	o-Xylene	0.832	0.860	-3.4	102	0.00
69 M	Styrene	1.331	1.376	-3.4	103	0.00
70	Isopropylbenzene	2.128	2.247	-5.6	105	0.00
71 M	1,1,2,2-Tetrachloroethane	0.735	0.757	-3.0	103	0.00
72	1,2,3-Trichloropropane	0.563	0.612	-8.7	104	0.00
73	Bromobenzene	0.496	0.519	-4.6	107	0.00
74	n-propylbenzene	2.412	2.469	-2.4	105	0.00
75	2-Chlorotoluene	1.508	1.565	-3.8	104	0.00
76	1,3,5-Trimethylbenzene	1.769	1.836	-3.8	105	0.00
77	t-1,4-Dichloro-2-butene	0.178	0.202	-13.5	125	0.00
78	4-Chlorotoluene	1.451	1.496	-3.1	105	0.00
79	tert-butylbenzene	1.564	1.655	-5.8	107	0.00
80	1,2,4-Trimethylbenzene	1.736	1.807	-4.1	104	0.00
81	sec-Butylbenzene	2.108	2.152	-2.1	104	0.00
82	p-Isopropyltoluene	1.712	1.763	-3.0	105	0.00
83 M	1,3-Dichlorobenzene	0.874	0.923	-5.6	107	0.00
84 M	1,4-Dichlorobenzene	0.873	0.901	-3.2	105	0.00
85	n-Butylbenzene	1.393	1.370	1.7	108	0.00
86 T	Hexachloroethane	0.335	0.338	-0.9	108	0.00
87 M	1,2-Dichlorobenzene	0.914	0.955	-4.5	106	0.00
88	1,2-Dibromo-3-Chloropropane	0.187	0.190	-1.6	103	0.00
89	1,2,4-Trichlorobenzene	0.480	0.479	0.2	108	0.00
90	Hexachlorobutadiene	0.194	0.205	-5.7	112	0.00
91 M	Naphthalene	1.874	1.980	-5.7	111	0.00
92	1,2,3-Trichlorobenzene	0.489	0.502	-2.7	110	0.00

(#) = Out of Range

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