

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121721\
 Data File : VN070155.D
 Acq On : 17 Dec 2021 10:58
 Operator : JC/MD
 Sample : VSTDCCC020
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Dec 18 05:28:58 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N121321W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Dec 14 00:59:04 2021
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	106	0.00
2 M	Dichlorodifluoromethane	2.177	2.142	1.6	103	0.00
3 M	Chloromethane	2.494	2.486	0.3	107	0.00
4 M	Vinyl Chloride	2.346	2.212	5.7	103	0.00
5 M	Bromomethane	1.283	1.285	-0.2	110	0.00
6 M	Chloroethane	1.347	1.291	4.2	104	0.00
7 M	Trichlorofluoromethane	3.063	2.774	9.4	97	0.00
8 T	Diethyl Ether	1.216	1.177	3.2	109	0.00
9	1,1,2-Trichlorotrifluoroeth	1.799	1.741	3.2	105	0.00
10 M	1,1-Dichloroethene	1.711	1.602	6.4	103	0.00
11	Methyl Iodide	2.395	2.156	10.0	100	0.00
12	Methyl Acetate	4.216	4.588	-8.8	126	0.00
13 M	Acrolein	0.341	0.342	-0.3	120	0.00
14 M	Acrylonitrile	1.217	1.257	-3.3	120	0.00
15 M	Acetone	0.442	0.511	-15.6	116	0.00
16 M	Carbon Disulfide	5.157	4.110	20.3	88	0.00
17	Allyl chloride	3.849	3.654	5.1	108	0.00
18 M	Methylene Chloride	2.063	1.972	4.4	107	0.00
19 M	trans-1,2-Dichloroethene	1.818	1.732	4.7	105	0.00
20 T	Diisopropyl ether	7.372	7.394	-0.3	113	0.00
21 M	1,1-Dichloroethane	3.886	3.725	4.1	108	0.00
22 M	cis-1,2-Dichloroethene	2.152	2.052	4.6	108	0.00
23 M	tert-Butyl Alcohol	0.385	0.441	-14.5	147	0.01
24 M	Methyl tert-Butyl Ether	6.077	6.319	-4.0	119	0.00
25 M	Chloroform	3.861	3.653	5.4	105	0.00
26	Cyclohexane	3.597	3.561	1.0	111	0.00
27 s	1,2-Dichloroethane-d4	2.659	2.744	-3.2	110	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	111	0.00
29	1,1-Dichloropropene	0.517	0.470	9.1	105	0.00
30 M	2-Butanone	0.355	0.383	-7.9	125	0.00
31	2,2-Dichloropropane	0.540	0.505	6.5	113	0.00
32 M	1,1,1-Trichloroethane	0.598	0.518	13.4	103	0.00
33 M	Carbon Tetrachloride	0.532	0.428	19.5	94	0.00
34 M	Benzene	1.527	1.407	7.9	107	0.00
35	Methacrylonitrile	0.329	0.317	3.6	122	0.00
36 M	1,2-Dichloroethane	0.603	0.554	8.1	108	0.00
37 M	Trichloroethene	0.363	0.322	11.3	105	0.00
38	Methylcyclohexane	0.614	0.564	8.1	109	0.00
39 M	1,2-Dichloropropane	0.407	0.377	7.4	109	0.00
40	Dibromomethane	0.269	0.238	11.5	103	0.00
41 M	Bromodichloromethane	0.568	0.456	19.7	94	0.00
42 M	Vinyl Acetate	1.044	1.009	3.4	118	0.00
43	Ethyl Acetate	0.617	0.629	-1.9	128	0.00
44	Isopropyl Acetate	0.994	0.977	1.7	121	0.00
45 T	1,4-Dioxane	0.007	0.008#	-14.3	137	0.00
46	Methyl methacrylate	0.471	0.454	3.6	117	0.00
47	n-amyl Acetate	0.693	0.659	4.9	133	0.00
48 M	t-1,3-Dichloropropene	0.571	0.480	15.9	106	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121721\
 Data File : VN070155.D
 Acq On : 17 Dec 2021 10:58
 Operator : JC/MD
 Sample : VSTDCCC020
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Dec 18 05:28:58 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N121321W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Dec 14 00:59:04 2021
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 30% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.620	0.533	14.0	106	0.00
50 M	1,1,2-Trichloroethane	0.365	0.338	7.4	109	0.00
51	Ethyl methacrylate	0.583	0.559	4.1	124	0.00
52	1,3-Dichloropropane	0.646	0.594	8.0	108	0.00
53 M	Dibromochloromethane	0.410	0.309	24.6	90	0.00
54 M	1,2-Dibromoethane	0.372	0.342	8.1	113	0.00
55 M	2-Chloroethyl vinyl ether	0.114	0.147	-28.9	176#	0.00
56 M	Bromoform	0.312	0.213	31.7#	85	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	112	0.00
58 M	4-Methyl-2-Pentanone	0.689	0.719	-4.4	123	0.00
59 M	2-Hexanone	0.509	0.564	-10.8	131	0.00
60 S	4-Bromofluorobenzene	0.494	0.489	1.0	113	0.00
61 M	Tetrachloroethene	0.350	0.328	6.3	104	0.00
62 M	Toluene	1.764	1.663	5.7	107	0.00
63 S	Toluene-d8	1.386	1.405	-1.4	110	0.00
64 M	Chlorobenzene	1.047	0.965	7.8	107	0.00
65	1,1,1,2-Tetrachloroethane	0.394	0.337	14.5	99	0.00
66 M	Ethyl Benzene	1.961	1.832	6.6	108	0.00
67 M	m/p-Xylenes	0.717	0.669	6.7	107	0.00
68 M	o-Xylene	0.709	0.666	6.1	109	0.00
69 M	Styrene	1.156	1.051	9.1	107	0.00
70	Isopropylbenzene	1.860	1.758	5.5	110	0.00
71 M	1,1,2,2-Tetrachloroethane	0.610	0.589	3.4	112	0.00
72	1,2,3-Trichloropropane	0.568	0.577	-1.6	112	0.00
73	Bromobenzene	0.432	0.394	8.8	107	0.00
74	n-propylbenzene	2.156	1.973	8.5	108	0.00
75	2-Chlorotoluene	1.319	1.224	7.2	109	0.00
76	1,3,5-Trimethylbenzene	1.528	1.412	7.6	108	0.00
77	t-1,4-Dichloro-2-butene	0.179	0.141	21.2	104	0.00
78	4-Chlorotoluene	1.265	1.152	8.9	108	0.00
79	tert-butylbenzene	1.326	1.231	7.2	111	0.00
80	1,2,4-Trimethylbenzene	1.478	1.375	7.0	108	0.00
81	sec-Butylbenzene	1.851	1.687	8.9	109	0.00
82	p-Isopropyltoluene	1.475	1.355	8.1	109	0.00
83 M	1,3-Dichlorobenzene	0.730	0.667	8.6	108	0.00
84 M	1,4-Dichlorobenzene	0.702	0.634	9.7	108	0.00
85	n-Butylbenzene	1.232	1.066	13.5	108	0.00
86 T	Hexachloroethane	0.294	0.207	29.6	86	0.00
87 M	1,2-Dichlorobenzene	0.701	0.660	5.8	109	0.00
88	1,2-Dibromo-3-Chloropropane	0.117	0.105	10.3	115	0.00
89	1,2,4-Trichlorobenzene	0.326	0.313	4.0	121	0.00
90	Hexachlorobutadiene	0.194	0.185	4.6	107	0.00
91 M	Naphthalene	0.847	0.909	-7.3	154#	0.00
92	1,2,3-Trichlorobenzene	0.308	0.315	-2.3	127	0.00

(#) = Out of Range

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