

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073649.D
 Acq On : 29 Jul 2022 21:04
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Aug 01 04:35:40 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N072722W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Jul 27 18:07:26 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	95	0.00
2 M	Dichlorodifluoromethane	2.039	1.998	2.0	92	0.00
3 M	Chloromethane	2.696	2.755	-2.2	92	0.00
4 M	Vinyl Chloride	3.309	3.396	-2.6	96	0.00
5 M	Bromomethane	1.681	1.861	-10.7	91	0.00
6 M	Chloroethane	2.242	2.278	-1.6	95	0.00
7 M	Trichlorofluoromethane	3.255	3.686	-13.2	109	0.00
8 T	Diethyl Ether	1.310	1.289	1.6	98	0.00
9	1,1,2-Trichlorotrifluoroeth	1.962	1.953	0.5	96	0.00
10 M	1,1-Dichloroethene	1.878	1.860	1.0	97	0.00
11	Methyl Iodide	2.166	2.077	4.1	100	0.00
12	Methyl Acetate	3.192	3.367	-5.5	101	0.00
13 M	Acrolein	0.561	0.631	-12.5	106	0.00
14 M	Acrylonitrile	1.473	1.496	-1.6	95	0.00
15 M	Acetone	0.410	0.432	-5.4	101	0.00
16 M	Carbon Disulfide	5.133	4.782	6.8	91	0.00
17	Allyl chloride	3.133	2.858	8.8	90	0.00
18 M	Methylene Chloride	2.380	2.585	-8.6	106	0.00
19 M	trans-1,2-Dichloroethene	2.082	2.042	1.9	95	-0.01
20 T	Diisopropyl ether	6.815	6.758	0.8	95	0.00
21 M	1,1-Dichloroethane	4.022	3.890	3.3	93	0.00
22 M	cis-1,2-Dichloroethene	2.505	2.516	-0.4	100	0.00
23 M	tert-Butyl Alcohol	0.521	0.537	-3.1	101	0.00
24 M	Methyl tert-Butyl Ether	6.619	6.734	-1.7	101	0.00
25 M	Chloroform	4.151	4.124	0.7	94	0.00
26	Cyclohexane	3.469	3.186	8.2	90	0.00
27 s	1,2-Dichloroethane-d4	2.594	2.699	-4.0	95	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	104	0.00
29	1,1-Dichloropropene	0.561	0.502	10.5	95	0.00
30 M	2-Butanone	0.387	0.367	5.2	98	0.00
31	2,2-Dichloropropane	0.585	0.455	22.2	82	0.00
32 M	1,1,1-Trichloroethane	0.649	0.618	4.8	99	0.00
33 M	Carbon Tetrachloride	0.556	0.530	4.7	102	0.00
34 M	Benzene	1.814	1.687	7.0	96	0.00
35	Methacrylonitrile	0.315	0.296	6.0	87	0.00
36 M	1,2-Dichloroethane	0.598	0.569	4.8	96	0.00
37 M	Trichloroethene	0.446	0.409	8.3	98	0.00
38	Methylcyclohexane	0.696	0.610	12.4	95	0.00
39 M	1,2-Dichloropropane	0.458	0.409	10.7	91	0.00
40	Dibromomethane	0.311	0.299	3.9	99	0.00
41 M	Bromodichloromethane	0.608	0.565	7.1	97	0.00
42 M	Vinyl Acetate	1.117	1.016	9.0	95	0.00
43	Ethyl Acetate	0.736	0.700	4.9	105	0.00
44	Isopropyl Acetate	1.023	0.954	6.7	98	0.00
45 T	1,4-Dioxane	0.012	0.011	8.3	97	0.00
46	Methyl methacrylate	0.457	0.403	11.8	94	0.00
47	n-amyl Acetate	0.816	0.740	9.3	101	0.00
48 M	t-1,3-Dichloropropene	0.626	0.553	11.7	100	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN072922\
 Data File : VN073649.D
 Acq On : 29 Jul 2022 21:04
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Aug 01 04:35:40 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N072722W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Wed Jul 27 18:07:26 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)
49 T	cis-1,3-Dichloropropene	0.695	0.623	10.4	97	0.00
50 M	1,1,2-Trichloroethane	0.466	0.446	4.3	101	0.00
51	Ethyl methacrylate	0.700	0.632	9.7	100	0.00
52	1,3-Dichloropropane	0.785	0.729	7.1	98	0.00
53 M	Dibromochloromethane	0.472	0.443	6.1	101	0.00
54 M	1,2-Dibromoethane	0.475	0.444	6.5	100	0.00
55 M	2-Chloroethyl vinyl ether	0.314	0.316	-0.6	108	0.00
56 M	Bromoform	0.371	0.334	10.0	105	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	103	0.00
58 M	4-Methyl-2-Pentanone	0.776	0.748	3.6	96	0.00
59 M	2-Hexanone	0.597	0.569	4.7	96	0.00
60 S	4-Bromofluorobenzene	0.531	0.521	1.9	103	0.00
61 M	Tetrachloroethene	0.460	0.395	14.1	88	0.00
62 M	Toluene	2.086	1.971	5.5	97	0.00
63 S	Toluene-d8	1.689	1.684	0.3	99	0.00
64 M	Chlorobenzene	1.246	1.162	6.7	99	0.00
65	1,1,1,2-Tetrachloroethane	0.463	0.432	6.7	99	0.00
66 M	Ethyl Benzene	2.230	2.006	10.0	95	0.00
67 M	m/p-Xylenes	0.852	0.806	5.4	99	0.00
68 M	o-Xylene	0.847	0.798	5.8	100	0.00
69 M	Styrene	1.396	1.286	7.9	97	0.00
70	Isopropylbenzene	2.165	2.015	6.9	98	0.00
71 M	1,1,2,2-Tetrachloroethane	0.761	0.755	0.8	102	0.00
72	1,2,3-Trichloropropane	0.636	0.654	-2.8	117	0.00
73	Bromobenzene	0.538	0.501	6.9	102	0.00
74	n-propylbenzene	2.519	2.270	9.9	96	0.00
75	2-Chlorotoluene	1.510	1.391	7.9	96	0.00
76	1,3,5-Trimethylbenzene	1.813	1.689	6.8	97	0.00
77	t-1,4-Dichloro-2-butene	0.232	0.200	13.8	98	0.00
78	4-Chlorotoluene	1.451	1.306	10.0	96	0.00
79	tert-butylbenzene	1.568	1.468	6.4	101	0.00
80	1,2,4-Trimethylbenzene	1.796	1.651	8.1	96	0.00
81	sec-Butylbenzene	2.251	2.016	10.4	97	0.00
82	p-Isopropyltoluene	1.848	1.642	11.1	98	0.00
83 M	1,3-Dichlorobenzene	0.984	0.905	8.0	101	0.00
84 M	1,4-Dichlorobenzene	0.949	0.858	9.6	98	0.00
85	n-Butylbenzene	1.498	1.213	19.0	94	0.00
86 T	Hexachloroethane	0.339	0.316	6.8	103	0.00
87 M	1,2-Dichlorobenzene	0.995	0.935	6.0	99	0.00
88	1,2-Dibromo-3-Chloropropane	0.166	0.157	5.4	101	0.00
89	1,2,4-Trichlorobenzene	0.548	0.448	18.2	99	0.00
90	Hexachlorobutadiene	0.277	0.252	9.0	103	0.00
91 M	Naphthalene	1.798	1.553	13.6	101	0.00
92	1,2,3-Trichlorobenzene	0.563	0.495	12.1	102	0.00

(#) = Out of Range

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