

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052622\
 Data File : VN072645.D
 Acq On : 26 May 2022 22:18
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: May 27 05:36:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N052622W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri May 27 05:07:24 2022
 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	103	0.00
2 M	Dichlorodifluoromethane	2.050	1.921	6.3	87	0.00
3 M	Chloromethane	2.749	2.343	14.8	90	0.00
4 M	Vinyl Chloride	2.298	1.993	13.3	89	0.00
5 M	Bromomethane	1.118	0.695	37.8#	69	0.02
6 M	Chloroethane	1.438	1.193	17.0	81	0.01
7 M	Trichlorofluoromethane	3.293	3.189	3.2	88	0.00
8 T	Diethyl Ether	1.382	1.369	0.9	87	0.00
9	1,1,2-Trichlorotrifluoroeth	1.840	1.859	-1.0	89	0.00
10 M	1,1-Dichloroethene	1.807	1.871	-3.5	91	0.00
11	Methyl Iodide	1.730	1.060	38.7#	56	0.00
12	Methyl Acetate	3.545	3.407	3.9	91	0.00
13 M	Acrolein	0.406	0.358	11.8	90	0.00
14 M	Acrylonitrile	1.527	1.480	3.1	89	0.00
15 M	Acetone	0.393	0.377	4.1	86	0.00
16 M	Carbon Disulfide	4.955	4.847	2.2	90	0.00
17	Allyl chloride	3.753	3.611	3.8	88	0.00
18 M	Methylene Chloride	2.375	2.356	0.8	86	0.00
19 M	trans-1,2-Dichloroethene	1.976	1.949	1.4	85	0.00
20 T	Diisopropyl ether	8.606	7.719	10.3	87	0.00
21 M	1,1-Dichloroethane	4.381	4.116	6.0	90	0.00
22 M	cis-1,2-Dichloroethene	2.510	2.391	4.7	87	0.00
23 M	tert-Butyl Alcohol	0.513	0.435	15.2	84	0.00
24 M	Methyl tert-Butyl Ether	6.941	6.845	1.4	90	0.00
25 M	Chloroform	4.451	4.163	6.5	92	0.00
26	Cyclohexane	4.132	3.582	13.3	88	0.00
27 s	1,2-Dichloroethane-d4	2.871	2.735	4.7	101	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	103	0.00
29	1,1-Dichloropropene	0.567	0.512	9.7	85	0.00
30 M	2-Butanone	0.382	0.366	4.2	87	0.00
31	2,2-Dichloropropane	0.575	0.509	11.5	80	0.00
32 M	1,1,1-Trichloroethane	0.629	0.600	4.6	88	0.00
33 M	Carbon Tetrachloride	0.559	0.513	8.2	87	0.00
34 M	Benzene	1.783	1.701	4.6	89	0.00
35	Methacrylonitrile	0.390	0.381	2.3	100	0.00
36 M	1,2-Dichloroethane	0.580	0.560	3.4	91	0.00
37 M	Trichloroethene	0.424	0.391	7.8	88	0.00
38	Methylcyclohexane	0.701	0.619	11.7	86	0.00
39 M	1,2-Dichloropropane	0.476	0.435	8.6	86	0.00
40	Dibromomethane	0.302	0.280	7.3	88	0.00
41 M	Bromodichloromethane	0.605	0.559	7.6	88	0.00
42 M	Vinyl Acetate	1.247	1.204	3.4	89	0.00
43	Ethyl Acetate	0.733	0.736	-0.4	96	0.00
44	Isopropyl Acetate	1.142	1.065	6.7	89	0.00
45 T	1,4-Dioxane	0.010	0.009#	10.0	86	0.00
46	Methyl methacrylate	0.538	0.463	13.9	87	0.00
47	n-amyl Acetate	0.832	0.731	12.1	87	0.00
48 M	t-1,3-Dichloropropene	0.653	0.570	12.7	86	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN052622\
 Data File : VN072645.D
 Acq On : 26 May 2022 22:18
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: May 27 05:36:24 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N052622W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri May 27 05:07:24 2022
 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	0.719	0.658	8.5	87	0.00
50 M	1,1,2-Trichloroethane	0.443	0.433	2.3	91	0.00
51	Ethyl methacrylate	0.748	0.670	10.4	86	0.00
52	1,3-Dichloropropane	0.782	0.741	5.2	89	0.00
53 M	Dibromochloromethane	0.456	0.419	8.1	87	0.00
54 M	1,2-Dibromoethane	0.455	0.425	6.6	87	0.00
55 M	2-Chloroethyl vinyl ether	0.327	0.342	-4.6	110	0.00
56 M	Bromoform	0.329	0.289	12.2	82	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	98	0.00
58 M	4-Methyl-2-Pentanone	0.867	0.825	4.8	89	0.00
59 M	2-Hexanone	0.618	0.587	5.0	89	0.00
60 S	4-Bromofluorobenzene	0.527	0.532	-0.9	99	0.00
61 M	Tetrachloroethene	0.397	0.391	1.5	88	0.00
62 M	Toluene	2.106	2.046	2.8	87	0.00
63 S	Toluene-d8	1.706	1.776	-4.1	101	0.00
64 M	Chlorobenzene	1.224	1.165	4.8	87	0.00
65	1,1,1,2-Tetrachloroethane	0.482	0.452	6.2	86	0.00
66 M	Ethyl Benzene	2.323	2.068	11.0	84	0.00
67 M	m/p-Xylenes	0.844	0.787	6.8	85	0.00
68 M	o-Xylene	0.840	0.790	6.0	86	0.00
69 M	Styrene	1.331	1.223	8.1	86	0.00
70	Isopropylbenzene	2.146	1.963	8.5	83	0.00
71 M	1,1,2,2-Tetrachloroethane	0.737	0.702	4.7	86	0.00
72	1,2,3-Trichloropropane	0.622	0.637	-2.4	90	0.00
73	Bromobenzene	0.472	0.457	3.2	88	0.00
74	n-propylbenzene	2.427	2.214	8.8	86	0.00
75	2-Chlorotoluene	1.472	1.371	6.9	87	0.00
76	1,3,5-Trimethylbenzene	1.694	1.564	7.7	86	0.00
77	t-1,4-Dichloro-2-butene	0.243	0.211	13.2	84	0.00
78	4-Chlorotoluene	1.340	1.260	6.0	88	0.00
79	tert-butylbenzene	1.482	1.413	4.7	86	0.00
80	1,2,4-Trimethylbenzene	1.635	1.924	-17.7	108	0.14
81	sec-Butylbenzene	2.144	1.924	10.3	86	0.00
82	p-Isopropyltoluene	1.662	1.465	11.9	85	0.00
83 M	1,3-Dichlorobenzene	0.818	0.733	10.4	87	0.00
84 M	1,4-Dichlorobenzene	0.795	0.717	9.8	87	0.00
85	n-Butylbenzene	1.366	1.160	15.1	87	0.00
86 T	Hexachloroethane	0.331	0.305	7.9	89	0.00
87 M	1,2-Dichlorobenzene	0.808	0.755	6.6	86	0.00
88	1,2-Dibromo-3-Chloropropane	0.143	0.130	9.1	89	0.00
89	1,2,4-Trichlorobenzene	0.399	0.345	13.5	85	0.00
90	Hexachlorobutadiene	0.200	0.180	10.0	82	0.00
91 M	Naphthalene	1.399	1.292	7.6	96	0.00
92	1,2,3-Trichlorobenzene	0.407	0.384	5.7	92	0.00

(#) = Out of Range

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