

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060723\
 Data File : VX036050.D
 Acq On : 07 Jun 2023 08:53
 Operator : JC/MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Jun 08 07:20:44 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X052323W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue May 23 11:36:37 2023
 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	101	0.00
2 M	Dichlorodifluoromethane	1.534	1.893	-23.4	194#	0.00
3 M	Chloromethane	1.882	1.811	3.8	110	0.00
4 M	Vinyl Chloride	1.769	1.784	-0.8	123	0.00
5 M	Bromomethane	1.190	1.137	4.5	105	0.00
6 M	Chloroethane	1.115	1.076	3.5	112	0.00
7 M	Trichlorofluoromethane	2.681	2.875	-7.2	152#	0.00
8 T	Diethyl Ether	1.115	1.069	4.1	102	0.00
9	1,1,2-Trichlorotrifluoroeth	1.443	1.670	-15.7	176#	0.00
10 M	1,1-Dichloroethene	1.504	1.492	0.8	119	0.00
11	Methyl Iodide	1.817	1.410	22.4	90	0.00
12	Methyl Acetate	4.249	4.637	-9.1	112	0.00
13 M	Acrolein	0.336	0.345	-2.7	114	0.00
14 M	Acrylonitrile	1.065	1.104	-3.7	109	0.00
15 M	Acetone	0.319	0.370	-16.0	117	-0.01
16 M	Carbon Disulfide	3.264	3.174	2.8	127	0.00
17	Allyl chloride	3.034	3.067	-1.1	114	0.00
18 M	Methylene Chloride	1.958	1.803	7.9	100	0.00
19 M	trans-1,2-Dichloroethene	1.695	1.636	3.5	106	0.00
20 T	Diisopropyl ether	6.550	6.509	0.6	106	0.00
21 M	1,1-Dichloroethane	3.433	3.508	-2.2	113	0.00
22 M	cis-1,2-Dichloroethene	2.098	2.016	3.9	106	0.00
23 M	tert-Butyl Alcohol	0.389	0.417	-7.2	110	-0.04
24 M	Methyl tert-Butyl Ether	6.061	6.242	-3.0	111	0.00
25 M	Chloroform	3.523	3.792	-7.6	120	0.00
26	Cyclohexane	2.596	2.675	-3.0	154#	0.00
27 s	1,2-Dichloroethane-d4	2.493	2.893	-16.0	120	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	104	0.00
29	1,1-Dichloropropene	0.418	0.445	-6.5	129	0.00
30 M	2-Butanone	0.273	0.312	-14.3	116	-0.02
31	2,2-Dichloropropane	0.449	0.483	-7.6	126	0.00
32 M	1,1,1-Trichloroethane	0.488	0.546	-11.9	132	0.00
33 M	Carbon Tetrachloride	0.385	0.442	-14.8	143	0.00
34 M	Benzene	1.318	1.328	-0.8	108	0.00
35	Methacrylonitrile	0.292	0.319	-9.2	114	0.00
36 M	1,2-Dichloroethane	0.513	0.598	-16.6	122	0.00
37 M	Trichloroethene	0.331	0.326	1.5	109	0.00
38	Methylcyclohexane	0.458	0.472	-3.1	163#	0.00
39 M	1,2-Dichloropropane	0.355	0.356	-0.3	108	0.00
40	Dibromomethane	0.240	0.255	-6.3	112	0.00
41 M	Bromodichloromethane	0.434	0.479	-10.4	120	0.00
42 M	Vinyl Acetate	0.803	0.830	-3.4	111	0.00
43	Ethyl Acetate	0.522	0.550	-5.4	113	0.00
44	Isopropyl Acetate	0.850	0.862	-1.4	109	0.00
45 T	1,4-Dioxane	0.008	0.009	-12.5	103	-0.04
46	Methyl methacrylate	0.411	0.439	-6.8	116	0.00
47	n-amyl Acetate	0.714	0.724	-1.4	110	0.00
48 M	t-1,3-Dichloropropene	0.480	0.497	-3.5	118	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.530	0.542	-2.3	111	0.00
50 M	1,1,2-Trichloroethane	0.343	0.361	-5.2	110	0.00
51	Ethyl methacrylate	0.529	0.541	-2.3	112	0.00
52	1,3-Dichloropropane	0.598	0.624	-4.3	108	0.00
53 M	Dibromochloromethane	0.315	0.330	-4.8	117	0.00
54 M	1,2-Dibromoethane	0.357	0.369	-3.4	109	0.00
55 M	2-Chloroethyl vinyl ether	0.281	0.295	-5.0	104	0.00
56 M	Bromoform	0.204	0.211	-3.4	124	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	106	0.00
58 M	4-Methyl-2-Pentanone	0.583	0.632	-8.4	114	0.00
59 M	2-Hexanone	0.438	0.488	-11.4	115	0.00
60 S	4-Bromofluorobenzene	0.503	0.556	-10.5	118	0.00
61 M	Tetrachloroethene	0.289	0.286	1.0	114	0.00
62 M	Toluene	1.545	1.552	-0.5	110	0.00
63 S	Toluene-d8	1.326	1.347	-1.6	108	0.00
64 M	Chlorobenzene	0.978	0.965	1.3	108	0.00
65	1,1,1,2-Tetrachloroethane	0.342	0.352	-2.9	116	0.00
66 M	Ethyl Benzene	1.728	1.777	-2.8	117	0.00
67 M	m/p-Xylenes	0.654	0.660	-0.9	113	0.00
68 M	o-Xylene	0.654	0.651	0.5	109	0.00
69 M	Styrene	1.080	1.099	-1.8	112	0.00
70	Isopropylbenzene	1.670	1.733	-3.8	119	0.00
71 M	1,1,2,2-Tetrachloroethane	0.591	0.617	-4.4	110	0.00
72	1,2,3-Trichloropropane	0.515	0.541	-5.0	117	0.00
73	Bromobenzene	0.406	0.400	1.5	109	0.00
74	n-propylbenzene	1.990	2.052	-3.1	119	0.00
75	2-Chlorotoluene	1.223	1.317	-7.7	122	0.00
76	1,3,5-Trimethylbenzene	1.418	1.511	-6.6	121	0.00
77	t-1,4-Dichloro-2-butene	0.155	0.143	7.7	116	0.00
78	4-Chlorotoluene	1.393	1.502	-7.8	121	0.00
79	tert-butylbenzene	1.369	1.394	-1.8	119	0.00
80	1,2,4-Trimethylbenzene	1.425	1.516	-6.4	120	0.00
81	sec-Butylbenzene	1.728	1.729	-0.1	122	0.00
82	p-Isopropyltoluene	1.438	1.445	-0.5	121	0.00
83 M	1,3-Dichlorobenzene	0.768	0.765	0.4	111	0.00
84 M	1,4-Dichlorobenzene	0.771	0.778	-0.9	112	0.00
85	n-Butylbenzene	1.279	1.281	-0.2	129	0.00
86 T	Hexachloroethane	0.226	0.217	4.0	129	0.00
87 M	1,2-Dichlorobenzene	0.752	0.763	-1.5	108	0.00
88	1,2-Dibromo-3-Chloropropane	0.113	0.129	-14.2	130	0.00
89	1,2,4-Trichlorobenzene	0.451	0.412	8.6	105	0.00
90	Hexachlorobutadiene	0.166	0.158	4.8	132	0.00
91 M	Naphthalene	1.560	1.480	5.1	104	0.00
92	1,2,3-Trichlorobenzene	0.450	0.407	9.6	103	0.00

(#) = Out of Range

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