

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX093022\
 Data File : VX031812.D
 Acq On : 30 Sep 2022 14:09
 Operator : JC/MD
 Sample : VSTDICV020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ICVVX093022

Quant Time: Oct 03 01:03:18 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X093022W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Mon Oct 03 00:56:55 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	102	0.00
2 M	Dichlorodifluoromethane	1.251	1.375	-9.9	110	0.00
3 M	Chloromethane	1.311	1.454	-10.9	104	0.00
4 M	Vinyl Chloride	1.758	1.902	-8.2	102	0.00
5 M	Bromomethane	1.609	1.787	-11.1	114	0.00
6 M	Chloroethane	1.550	1.494	3.6	96	0.00
7 M	Trichlorofluoromethane	3.758	4.109	-9.3	103	0.00
8 T	Diethyl Ether	1.273	1.421	-11.6	103	0.00
9	1,1,2-Trichlorotrifluoroeth	1.783	1.980	-11.0	107	0.00
10 M	1,1-Dichloroethene	1.647	1.720	-4.4	103	0.00
11	Methyl Iodide	1.803	1.410	21.8	92	0.00
12	Methyl Acetate	1.882	2.054	-9.1	104	0.00
13 M	Acrolein	0.167	0.163	2.4	106	0.00
14 M	Acrylonitrile	0.799	0.843	-5.5	100	0.00
15 M	Acetone	0.216	0.256	-18.5	109	0.00
16 M	Carbon Disulfide	3.613	3.721	-3.0	107	0.00
17	Allyl chloride	1.994	2.116	-6.1	103	0.00
18 M	Methylene Chloride	1.885	2.021	-7.2	103	0.00
19 M	trans-1,2-Dichloroethene	1.907	1.996	-4.7	103	0.00
20 T	Diisopropyl ether	4.582	4.992	-8.9	102	0.00
21 M	1,1-Dichloroethane	2.989	3.181	-6.4	103	0.00
22 M	cis-1,2-Dichloroethene	2.304	2.444	-6.1	106	0.00
23 M	tert-Butyl Alcohol	0.298	0.329	-10.4	102	-0.01
24 M	Methyl tert-Butyl Ether	5.661	6.003	-6.0	103	0.00
25 M	Chloroform	3.670	3.935	-7.2	103	0.00
26	Cyclohexane	2.466	2.697	-9.4	104	0.00
27 s	1,2-Dichloroethane-d4	2.258	2.369	-4.9	100	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	104	0.00
29	1,1-Dichloropropene	0.465	0.489	-5.2	106	0.00
30 M	2-Butanone	0.189	0.203	-7.4	101	0.00
31	2,2-Dichloropropane	0.466	0.476	-2.1	107	0.00
32 M	1,1,1-Trichloroethane	0.616	0.624	-1.3	105	0.00
33 M	Carbon Tetrachloride	0.564	0.573	-1.6	105	0.00
34 M	Benzene	1.376	1.439	-4.6	104	0.00
35	Methacrylonitrile	0.198	0.207	-4.5	102	0.00
36 M	1,2-Dichloroethane	0.485	0.521	-7.4	104	0.00
37 M	Trichloroethene	0.449	0.455	-1.3	107	0.00
38	Methylcyclohexane	0.569	0.594	-4.4	107	0.00
39 M	1,2-Dichloropropane	0.315	0.333	-5.7	107	0.00
40	Dibromomethane	0.275	0.281	-2.2	101	0.00
41 M	Bromodichloromethane	0.499	0.513	-2.8	106	0.00
42 M	Vinyl Acetate	0.683	0.725	-6.1	104	0.00
43	Ethyl Acetate	0.380	0.390	-2.6	98	0.00
44	Isopropyl Acetate	0.605	0.628	-3.8	103	0.00
45 T	1,4-Dioxane	0.009	0.009	0.0	94	0.00
46	Methyl methacrylate	0.282	0.299	-6.0	102	0.00
47	n-amyl Acetate	0.519	0.548	-5.6	106	0.00
48 M	t-1,3-Dichloropropene	0.492	0.493	-0.2	106	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	0.540	0.535	0.9	102	0.00
50 M	1,1,2-Trichloroethane	0.398	0.422	-6.0	106	0.00
51	Ethyl methacrylate	0.517	0.537	-3.9	103	0.00
52	1,3-Dichloropropane	0.611	0.650	-6.4	104	0.00
53 M	Dibromochloromethane	0.444	0.445	-0.2	108	0.00
54 M	1,2-Dibromoethane	0.441	0.458	-3.9	105	0.00
55 M	2-Chloroethyl vinyl ether	0.253	0.243	4.0	91	0.00
56 M	Bromoform	0.325	0.303	6.8	106	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	101	0.00
58 M	4-Methyl-2-Pentanone	0.403	0.442	-9.7	101	0.00
59 M	2-Hexanone	0.302	0.333	-10.3	102	0.00
60 S	4-Bromofluorobenzene	0.555	0.578	-4.1	102	0.00
61 M	Tetrachloroethene	0.482	0.491	-1.9	105	0.00
62 M	Toluene	1.718	1.855	-8.0	107	0.00
63 S	Toluene-d8	1.524	1.579	-3.6	102	0.00
64 M	Chlorobenzene	1.172	1.240	-5.8	105	0.00
65	1,1,1,2-Tetrachloroethane	0.444	0.467	-5.2	109	0.00
66 M	Ethyl Benzene	1.936	2.065	-6.7	104	0.00
67 M	m/p-Xylenes	0.806	0.866	-7.4	104	0.00
68 M	o-Xylene	0.791	0.862	-9.0	107	0.00
69 M	Styrene	1.285	1.384	-7.7	105	0.00
70	Isopropylbenzene	2.057	2.254	-9.6	106	0.00
71 M	1,1,2,2-Tetrachloroethane	0.586	0.643	-9.7	105	0.00
72	1,2,3-Trichloropropane	0.508	0.558	-9.8	103	0.00
73	Bromobenzene	0.552	0.588	-6.5	108	0.00
74	n-propylbenzene	2.261	2.482	-9.8	106	0.00
75	2-Chlorotoluene	1.359	1.501	-10.4	106	0.00
76	1,3,5-Trimethylbenzene	1.734	1.918	-10.6	106	0.00
77	t-1,4-Dichloro-2-butene	0.147	0.136	7.5	104	0.00
78	4-Chlorotoluene	1.550	1.729	-11.5	107	0.00
79	tert-butylbenzene	1.743	1.888	-8.3	105	0.00
80	1,2,4-Trimethylbenzene	1.719	1.904	-10.8	106	0.00
81	sec-Butylbenzene	2.134	2.344	-9.8	107	0.00
82	p-Isopropyltoluene	1.876	2.097	-11.8	108	0.00
83 M	1,3-Dichlorobenzene	1.044	1.136	-8.8	110	0.00
84 M	1,4-Dichlorobenzene	1.070	1.145	-7.0	107	0.00
85	n-Butylbenzene	1.446	1.603	-10.9	110	0.00
86 T	Hexachloroethane	0.294	0.298	-1.4	107	0.00
87 M	1,2-Dichlorobenzene	1.041	1.118	-7.4	108	0.00
88	1,2-Dibromo-3-Chloropropane	0.118	0.121	-2.5	101	0.00
89	1,2,4-Trichlorobenzene	0.674	0.699	-3.7	113	0.00
90	Hexachlorobutadiene	0.271	0.285	-5.2	117	0.00
91 M	Naphthalene	2.156	2.294	-6.4	105	0.00
92	1,2,3-Trichlorobenzene	0.667	0.713	-6.9	115	0.00

(#) = Out of Range

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