

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030924\
 Data File : VX040620.D
 Acq On : 08 Mar 2024 10:12
 Operator : JC/MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Mar 08 23:49:51 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X030524W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Mar 05 13:21:29 2024
 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	98	0.00
2 M	Dichlorodifluoromethane	2.035	2.016	0.9	109	0.00
3 M	Chloromethane	2.085	2.043	2.0	105	0.00
4 M	Vinyl Chloride	2.253	2.216	1.6	108	0.00
5 M	Bromomethane	1.376	1.878	-36.5#	126	0.01
6 M	Chloroethane	1.324	1.409	-6.4	126	0.02
7 M	Trichlorofluoromethane	3.430	3.398	0.9	115	0.01
8 T	Diethyl Ether	1.165	1.029	11.7	102	0.00
9	1,1,2-Trichlorotrifluoroeth	1.879	1.930	-2.7	113	0.00
10 M	1,1-Dichloroethene	1.789	1.813	-1.3	113	0.00
11	Methyl Iodide	2.164	1.489	31.2#	79	0.00
12	Methyl Acetate	2.947	2.811	4.6	106	0.00
13 M	Acrolein	0.556	0.443	20.3	90	0.00
14 M	Acrylonitrile	1.160	1.105	4.7	104	0.00
15 M	Acetone	0.332	0.324	2.4	100	0.00
16 M	Carbon Disulfide	4.917	4.585	6.8	106	0.00
17	Allyl chloride	3.301	3.262	1.2	110	0.00
18 M	Methylene Chloride	2.043	2.001	2.1	108	0.00
19 M	trans-1,2-Dichloroethene	1.953	1.887	3.4	106	0.00
20 T	Diisopropyl ether	6.603	6.548	0.8	110	0.00
21 M	1,1-Dichloroethane	3.799	3.775	0.6	109	0.00
22 M	cis-1,2-Dichloroethene	2.268	2.190	3.4	105	0.00
23 M	tert-Butyl Alcohol	0.548	0.484	11.7	101	-0.04
24 M	Methyl tert-Butyl Ether	6.602	6.414	2.8	108	0.00
25 M	Chloroform	4.158	4.182	-0.6	112	0.00
26	Cyclohexane	3.066	2.988	2.5	111	0.00
27 s	1,2-Dichloroethane-d4	2.996	2.887	3.6	101	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	103	0.00
29	1,1-Dichloropropene	0.519	0.531	-2.3	114	0.00
30 M	2-Butanone	0.317	0.315	0.6	104	0.00
31	2,2-Dichloropropane	0.560	0.578	-3.2	111	0.00
32 M	1,1,1-Trichloroethane	0.654	0.664	-1.5	112	0.00
33 M	Carbon Tetrachloride	0.551	0.548	0.5	111	0.00
34 M	Benzene	1.402	1.420	-1.3	109	0.00
35	Methacrylonitrile	0.325	0.322	0.9	105	0.00
36 M	1,2-Dichloroethane	0.648	0.659	-1.7	108	0.00
37 M	Trichloroethene	0.377	0.385	-2.1	112	0.00
38	Methylcyclohexane	0.559	0.590	-5.5	114	0.00
39 M	1,2-Dichloropropane	0.357	0.374	-4.8	111	0.00
40	Dibromomethane	0.280	0.299	-6.8	110	0.00
41 M	Bromodichloromethane	0.558	0.590	-5.7	113	0.00
42 M	Vinyl Acetate	1.091	1.082	0.8	108	0.00
43	Ethyl Acetate	0.597	0.571	4.4	99	0.00
44	Isopropyl Acetate	0.953	0.917	3.8	105	0.00
45 T	1,4-Dioxane	0.010	0.009	10.0	97	0.00
46	Methyl methacrylate	0.473	0.475	-0.4	105	0.00
47	n-amyl Acetate	0.821	0.696	15.2	102	0.00
48 M	t-1,3-Dichloropropene	0.590	0.565	4.2	108	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX030924\
 Data File : VX040620.D
 Acq On : 08 Mar 2024 10:12
 Operator : JC/MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Mar 08 23:49:51 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\624X030524W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Mar 05 13:21:29 2024
 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	0.601	0.621	-3.3	114	0.00
50 M	1,1,2-Trichloroethane	0.372	0.396	-6.5	113	0.00
51	Ethyl methacrylate	0.586	0.595	-1.5	109	0.00
52	1,3-Dichloropropane	0.636	0.657	-3.3	109	0.00
53 M	Dibromochloromethane	0.425	0.433	-1.9	111	0.00
54 M	1,2-Dibromoethane	0.412	0.416	-1.0	108	0.00
55 M	2-Chloroethyl vinyl ether	0.299	0.268	10.4	92	0.00
56 M	Bromoform	0.301	0.279	7.3	108	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	112	0.00
58 M	4-Methyl-2-Pentanone	0.668	0.669	-0.1	106	0.00
59 M	2-Hexanone	0.540	0.535	0.9	107	0.00
60 S	4-Bromofluorobenzene	0.515	0.452	12.2	100	0.00
61 M	Tetrachloroethene	0.356	0.376	-5.6	113	0.00
62 M	Toluene	1.652	1.689	-2.2	109	0.00
63 S	Toluene-d8	1.326	1.341	-1.1	101	0.00
64 M	Chlorobenzene	1.010	1.047	-3.7	124	0.00
65	1,1,1,2-Tetrachloroethane	0.385	0.397	-3.1	125	0.00
66 M	Ethyl Benzene	1.905	1.959	-2.8	124	0.00
67 M	m/p-Xylenes	0.704	0.698	0.9	120	0.00
68 M	o-Xylene	0.677	0.644	4.9	113	0.00
69 M	Styrene	1.138	1.070	6.0	111	0.00
70	Isopropylbenzene	1.830	1.788	2.3	117	0.00
71 M	1,1,2,2-Tetrachloroethane	0.642	0.589	8.3	108	0.00
72	1,2,3-Trichloropropane	0.577	0.511	11.4	111	0.00
73	Bromobenzene	0.433	0.411	5.1	113	0.00
74	n-propylbenzene	2.260	2.072	8.3	112	0.00
75	2-Chlorotoluene	1.341	1.215	9.4	111	0.00
76	1,3,5-Trimethylbenzene	1.611	1.514	6.0	115	0.00
77	t-1,4-Dichloro-2-butene	0.179	0.148	17.3	100	0.00
78	4-Chlorotoluene	1.600	1.485	7.2	118	0.00
79	tert-butylbenzene	1.504	1.358	9.7	114	0.00
80	1,2,4-Trimethylbenzene	1.587	1.432	9.8	114	0.00
81	sec-Butylbenzene	1.985	1.818	8.4	115	0.00
82	p-Isopropyltoluene	1.625	1.488	8.4	115	0.00
83 M	1,3-Dichlorobenzene	0.843	0.780	7.5	114	0.00
84 M	1,4-Dichlorobenzene	0.862	0.789	8.5	113	0.00
85	n-Butylbenzene	1.581	1.604	-1.5	128	0.00
86 T	Hexachloroethane	0.284	0.277	2.5	128	0.00
87 M	1,2-Dichlorobenzene	0.819	0.838	-2.3	124	0.00
88	1,2-Dibromo-3-Chloropropane	0.172	0.167	2.9	119	0.00
89	1,2,4-Trichlorobenzene	0.552	0.566	-2.5	126	0.00
90	Hexachlorobutadiene	0.223	0.235	-5.4	130	0.00
91 M	Naphthalene	1.799	1.744	3.1	120	0.00
92	1,2,3-Trichlorobenzene	0.544	0.555	-2.0	125	0.00

(#) = Out of Range

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