

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX091120\
 Data File : VX018373.D
 Acq On : 11 Sep 2020 20:01
 Operator : JC/SP
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Sep 12 07:25:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X090420W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Fri Sep 04 12:13:41 2020
 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	1.000	1.000	0.0	93	0.00
2 M	Dichlorodifluoromethane	1.587	1.501	5.4	94	0.00
3 M	Chloromethane	1.686	1.628	3.4	96	0.00
4 M	Vinyl Chloride	1.776	1.768	0.5	100	0.00
5 M	Bromomethane	1.208	1.252	-3.6	94	0.00
6 M	Chloroethane	1.079	1.057	2.0	96	0.00
7 M	Trichlorofluoromethane	3.009	3.058	-1.6	102	0.00
8 T	Diethyl Ether	0.949	0.945	0.4	99	0.00
9	1,1,2-Trichlorotrifluoroeth	1.537	1.468	4.5	94	0.00
10 M	1,1-Dichloroethene	1.509	1.442	4.4	95	0.00
11	Methyl Iodide	1.623	1.448	10.8	110	0.00
12	Methyl Acetate	1.987	2.044	-2.9	100	0.00
13 M	Acrolein	0.160	0.172	-7.5	120	0.00
14 M	Acrylonitrile	0.900	0.880	2.2	98	0.00
15 M	Acetone	0.245	0.236	3.7	95	0.00
16 M	Carbon Disulfide	4.394	4.175	5.0	94	0.00
17	Allyl chloride	2.697	2.652	1.7	100	0.00
18 M	Methylene Chloride	1.771	1.710	3.4	94	0.00
19 M	trans-1,2-Dichloroethene	1.676	1.654	1.3	100	0.00
20 T	Diisopropyl ether	5.340	5.243	1.8	96	0.00
21 M	1,1-Dichloroethane	3.065	2.993	2.3	96	0.00
22 M	cis-1,2-Dichloroethene	1.854	1.803	2.8	96	0.00
23 M	tert-Butyl Alcohol	0.390	0.385	1.3	103	0.00
24 M	Methyl tert-Butyl Ether	5.365	5.209	2.9	98	0.00
25 M	Chloroform	3.271	3.233	1.2	97	0.00
26	Cyclohexane	2.474	2.397	3.1	99	0.00
27 s	1,2-Dichloroethane-d4	2.367	2.366	0.0	91	0.00
28 I	1,4-Difluorobenzene	1.000	1.000	0.0	91	0.00
29	1,1-Dichloropropene	0.458	0.441	3.7	97	0.00
30 M	2-Butanone	0.251	0.245	2.4	96	0.00
31	2,2-Dichloropropane	0.553	0.490	11.4	88	0.00
32 M	1,1,1-Trichloroethane	0.599	0.600	-0.2	98	0.00
33 M	Carbon Tetrachloride	0.559	0.554	0.9	97	0.00
34 M	Benzene	1.307	1.299	0.6	97	0.00
35	Methacrylonitrile	0.264	0.264	0.0	100	0.00
36 M	1,2-Dichloroethane	0.541	0.541	0.0	99	0.00
37 M	Trichloroethene	0.378	0.365	3.4	98	0.00
38	Methylcyclohexane	0.488	0.454	7.0	99	0.00
39 M	1,2-Dichloropropane	0.350	0.337	3.7	94	0.00
40	Dibromomethane	0.253	0.248	2.0	96	0.00
41 M	Bromodichloromethane	0.524	0.513	2.1	97	0.00
42 M	Vinyl Acetate	0.935	0.935	0.0	99	0.00
43	Ethyl Acetate	0.499	0.485	2.8	96	0.00
44	Isopropyl Acetate	0.804	0.796	1.0	99	0.00
45 T	1,4-Dioxane	0.007	0.006#	14.3	91	0.00

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Quant Time: Sep 12 07:25:00 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X090420W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
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 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
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	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
46	Methyl methacrylate	0.395	0.386	2.3	103	0.00
47	n-amyl Acetate	0.706	0.680	3.7	101	0.00
48 M	t-1,3-Dichloropropene	0.563	0.535	5.0	96	0.00
49 T	cis-1,3-Dichloropropene	0.591	0.576	2.5	96	0.00
50 M	1,1,2-Trichloroethane	0.352	0.350	0.6	97	0.00
51	Ethyl methacrylate	0.510	0.489	4.1	99	0.00
52	1,3-Dichloropropane	0.585	0.565	3.4	96	0.00
53 M	Dibromochloromethane	0.438	0.432	1.4	101	0.00
54 M	1,2-Dibromoethane	0.386	0.378	2.1	98	0.00
55 M	2-Chloroethyl vinyl ether	0.267	0.273	-2.2	103	0.00
56 M	Bromoform	0.333	0.326	2.1	104	0.00
57 I	Chlorobenzene-d5	1.000	1.000	0.0	91	0.00
58 M	4-Methyl-2-Pentanone	0.516	0.524	-1.6	99	0.00
59 M	2-Hexanone	0.391	0.392	-0.3	99	0.00
60 S	4-Bromofluorobenzene	0.520	0.507	2.5	89	0.00
61 M	Tetrachloroethene	0.374	0.358	4.3	93	0.00
62 M	Toluene	1.510	1.481	1.9	97	0.00
63 S	Toluene-d8	1.322	1.317	0.4	89	0.00
64 M	Chlorobenzene	0.970	0.962	0.8	97	0.00
65	1,1,1,2-Tetrachloroethane	0.393	0.385	2.0	98	0.00
66 M	Ethyl Benzene	1.668	1.656	0.7	99	0.00
67 M	m/p-Xylenes	0.638	0.640	-0.3	99	0.00
68 M	o-Xylene	0.603	0.592	1.8	101	0.00
69 M	Styrene	1.052	1.027	2.4	98	0.00
70	Isopropylbenzene	1.676	1.628	2.9	97	0.00
71 M	1,1,2,2-Tetrachloroethane	0.548	0.537	2.0	97	0.00
72	1,2,3-Trichloropropane	0.526	0.519	1.3	96	0.00
73	Bromobenzene	0.459	0.448	2.4	97	0.00
74	n-propylbenzene	1.974	1.869	5.3	95	0.00
75	2-Chlorotoluene	1.196	1.197	-0.1	99	0.00
76	1,3,5-Trimethylbenzene	1.460	1.408	3.6	97	0.00
77	t-1,4-Dichloro-2-butene	0.205	0.192	6.3	98	0.00
78	4-Chlorotoluene	1.427	1.375	3.6	96	0.00
79	tert-butylbenzene	1.407	1.357	3.6	97	0.00
80	1,2,4-Trimethylbenzene	1.444	1.443	0.1	99	0.00
81	sec-Butylbenzene	1.643	1.588	3.3	99	0.00
82	p-Isopropyltoluene	1.532	1.457	4.9	97	0.00
83 M	1,3-Dichlorobenzene	0.821	0.793	3.4	95	0.00
84 M	1,4-Dichlorobenzene	0.821	0.813	1.0	97	0.00
85	n-Butylbenzene	1.349	1.244	7.8	96	0.00
86 T	Hexachloroethane	0.302	0.289	4.3	99	0.00
87 M	1,2-Dichlorobenzene	0.769	0.765	0.5	99	0.00
88	1,2-Dibromo-3-Chloropropane	0.140	0.137	2.1	103	0.00
89	1,2,4-Trichlorobenzene	0.520	0.482	7.3	96	0.00
90	Hexachlorobutadiene	0.220	0.211	4.1	93	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
91 M	Naphthalene	1.624	1.558	4.1	100	0.00
92	1,2,3-Trichlorobenzene	0.515	0.502	2.5	101	0.00

(#) = Out of Range

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