

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\WX091820\
 Data File : VX018506.D
 Acq On : 18 Sep 2020 20:27
 Operator : JC/SP
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Sep 21 03:58:05 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	30.000	30.000	0.0	119	0.01
2 M	Dichlorodifluoromethane	20.000	15.202	24.0	97	0.00
3 M	Chloromethane	20.000	15.272	23.6	98	0.00
4 M	Vinyl Chloride	20.000	16.402	18.0	105	0.00
5 M	Bromomethane	20.000	12.590	37.0#	73	0.00
6 M	Chloroethane	20.000	17.180	14.1	108	0.00
7 M	Trichlorofluoromethane	20.000	15.410	22.9	100	0.00
8 T	Diethyl Ether	20.000	16.461	17.7	105	0.00
9	1,1,2-Trichlorotrifluoroeth	20.000	15.851	20.7	101	0.00
10 M	1,1-Dichloroethene	20.000	16.739	16.3	107	0.00
11	Methyl Iodide	20.000	15.403	23.0	87	0.00
12	Methyl Acetate	20.000	18.219	8.9	114	0.00
13 M	Acrolein	100.000	68.565	31.4#	99	0.00
14 M	Acrylonitrile	100.000	82.395	17.6	106	0.00
15 M	Acetone	100.000	80.195	19.8	101	0.00
16 M	Carbon Disulfide	20.000	16.149	19.3	102	0.00
17	Allyl chloride	20.000	16.534	17.3	107	0.00
18 M	Methylene Chloride	20.000	15.864	20.7	99	0.00
19 M	trans-1,2-Dichloroethene	20.000	16.311	18.4	106	0.00
20 T	Diisopropyl ether	20.000	17.402	13.0	109	0.00
21 M	1,1-Dichloroethane	20.000	16.283	18.6	103	0.00
22 M	cis-1,2-Dichloroethene	20.000	16.882	15.6	107	0.00
23 M	tert-Butyl Alcohol	100.000	88.554	11.4	118	0.00
24 M	Methyl tert-Butyl Ether	20.000	16.296	18.5	105	0.00
25 M	Chloroform	20.000	16.067	19.7	101	0.00
26	Cyclohexane	20.000	16.481	17.6	108	0.00
27 s	1,2-Dichloroethane-d4	30.000	29.585	1.4	115	0.00
28 I	1,4-Difluorobenzene	30.000	30.000	0.0	121	0.00
29	1,1-Dichloropropene	20.000	15.943	20.3	106	0.00
30 M	2-Butanone	100.000	81.306	18.7	106	0.00
31	2,2-Dichloropropane	20.000	13.953	30.2#	92	0.00
32 M	1,1,1-Trichloroethane	20.000	15.784	21.1	103	0.00
33 M	Carbon Tetrachloride	20.000	15.410	22.9	100	0.00
34 M	Benzene	20.000	15.966	20.2	104	0.00
35	Methacrylonitrile	20.000	16.645	16.8	111	0.00
36 M	1,2-Dichloroethane	20.000	15.958	20.2	105	0.00
37 M	Trichloroethene	20.000	16.330	18.4	110	0.00
38	Methylcyclohexane	20.000	15.969	20.2	113	0.00
39 M	1,2-Dichloropropane	20.000	15.947	20.3	103	0.00
40	Dibromomethane	20.000	15.924	20.4	104	0.00
41 M	Bromodichloromethane	20.000	15.964	20.2	105	0.00
42 M	Vinyl Acetate	100.000	81.561	18.4	107	0.00
43	Ethyl Acetate	20.000	15.766	21.2	103	0.00
44	Isopropyl Acetate	20.000	16.276	18.6	108	0.00
45 T	1,4-Dioxane	400.000	292.428	26.9	100	0.00

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Quant Time: Sep 21 03:58:05 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	Amount	Calc.	%Dev	Area%	Dev(min)
46	Methyl methacrylate	20.000	16.042	19.8	112	0.00
47	n-amyl Acetate	20.000	15.751	21.2	110	0.00
48 M	t-1,3-Dichloropropene	20.000	15.763	21.2	106	0.00
49 T	cis-1,3-Dichloropropene	20.000	16.119	19.4	106	0.00
50 M	1,1,2-Trichloroethane	20.000	15.594	22.0	102	0.00
51	Ethyl methacrylate	20.000	16.207	19.0	112	0.00
52	1,3-Dichloropropane	20.000	16.179	19.1	106	0.00
53 M	Dibromochloromethane	20.000	15.453	22.7	105	0.00
54 M	1,2-Dibromoethane	20.000	15.739	21.3	105	0.00
55 M	2-Chloroethyl vinyl ether	100.000	84.456	15.5	113	0.00
56 M	Bromoform	20.000	15.715	21.4	110	0.00
57 I	Chlorobenzene-d5	30.000	30.000	0.0	121	0.00
58 M	4-Methyl-2-Pentanone	100.000	83.222	16.8	107	0.00
59 M	2-Hexanone	100.000	82.216	17.8	107	0.00
60 S	4-Bromofluorobenzene	30.000	29.236	2.5	118	0.00
61 M	Tetrachloroethene	20.000	15.440	22.8	99	0.00
62 M	Toluene	20.000	16.353	18.2	107	0.00
63 S	Toluene-d8	30.000	29.368	2.1	116	0.00
64 M	Chlorobenzene	20.000	16.555	17.2	107	0.00
65	1,1,1,2-Tetrachloroethane	20.000	15.787	21.1	104	0.00
66 M	Ethyl Benzene	20.000	16.146	19.3	107	0.00
67 M	m/p-Xylenes	40.000	32.579	18.6	107	0.00
68 M	o-Xylene	20.000	16.252	18.7	110	0.00
69 M	Styrene	20.000	16.164	19.2	108	0.00
70	Isopropylbenzene	20.000	15.934	20.3	105	0.00
71 M	1,1,2,2-Tetrachloroethane	20.000	16.519	17.4	108	0.00
72	1,2,3-Trichloropropane	20.000	15.948	20.3	103	0.00
73	Bromobenzene	20.000	16.478	17.6	108	0.00
74	n-propylbenzene	20.000	15.886	20.6	105	0.00
75	2-Chlorotoluene	20.000	15.537	22.3	102	0.00
76	1,3,5-Trimethylbenzene	20.000	15.799	21.0	105	0.00
77	t-1,4-Dichloro-2-butene	20.000	13.958	30.2#	96	0.00
78	4-Chlorotoluene	20.000	15.853	20.7	105	0.00
79	tert-butylbenzene	20.000	16.060	19.7	107	0.00
80	1,2,4-Trimethylbenzene	20.000	16.053	19.7	106	0.00
81	sec-Butylbenzene	20.000	15.605	22.0	105	0.00
82	p-Isopropyltoluene	20.000	15.560	22.2	105	0.00
83 M	1,3-Dichlorobenzene	20.000	15.887	20.6	104	0.00
84 M	1,4-Dichlorobenzene	20.000	16.360	18.2	106	0.00
85	n-Butylbenzene	20.000	15.481	22.6	107	0.00
86 T	Hexachloroethane	20.000	14.745	26.3	100	0.00
87 M	1,2-Dichlorobenzene	20.000	16.501	17.5	109	0.00
88	1,2-Dibromo-3-Chloropropane	20.000	15.399	23.0	107	0.00
89	1,2,4-Trichlorobenzene	20.000	16.168	19.2	111	0.00
90	Hexachlorobutadiene	20.000	15.902	20.5	102	0.00

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Response via : Initial Calibration

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
91 M	Naphthalene	20.000	16.535	17.3	114	0.00
92	1,2,3-Trichlorobenzene	20.000	16.691	16.5	114	0.00

(#) = Out of Range

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