

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX111219\
 Data File : VX013374.D
 Acq On : 12 Nov 2019 09:58
 Operator : JC/SP
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Nov 12 12:36:20 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X111119W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Nov 12 01:16:13 2019
 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	107	0.00
2 M	Dichlorodifluoromethane	20.000	18.003	10.0	100	0.00
3 M	Chloromethane	20.000	18.666	6.7	104	0.00
4 M	Vinyl Chloride	20.000	17.824	10.9	100	0.00
5 M	Bromomethane	20.000	17.603	12.0	111	0.00
6 M	Chloroethane	20.000	18.409	8.0	106	0.00
7 M	Trichlorofluoromethane	20.000	16.928	15.4	98	0.00
8 T	Diethyl Ether	20.000	17.565	12.2	102	0.00
9	1,1,2-Trichlorotrifluoroeth	20.000	17.487	12.6	100	0.00
10 M	1,1-Dichloroethene	20.000	17.458	12.7	101	0.00
11	Methyl Iodide	20.000	14.788	26.1	90	0.00
12	Methyl Acetate	20.000	17.316	13.4	99	0.00
13 M	Acrolein	100.000	89.592	10.4	103	0.00
14 M	Acrylonitrile	100.000	84.326	15.7	97	0.00
15 M	Acetone	100.000	99.572	0.4	101	0.00
16 M	Carbon Disulfide	20.000	17.371	13.1	102	0.00
17	Allyl chloride	20.000	17.564	12.2	101	0.00
18 M	Methylene Chloride	20.000	17.501	12.5	103	0.00
19 M	trans-1,2-Dichloroethene	20.000	17.896	10.5	105	0.00
20 T	Diisopropyl ether	20.000	17.860	10.7	103	0.00
21 M	1,1-Dichloroethane	20.000	17.799	11.0	103	0.00
22 M	cis-1,2-Dichloroethene	20.000	17.743	11.3	104	0.00
23 M	tert-Butyl Alcohol	100.000	83.125	16.9	101	0.00
24 M	Methyl tert-Butyl Ether	20.000	18.068	9.7	104	0.00
25 M	Chloroform	20.000	17.966	10.2	104	0.00
26	Cyclohexane	20.000	17.047	14.8	100	0.00
27 s	1,2-Dichloroethane-d4	30.000	28.522	4.9	105	0.00
28 I	1,4-Difluorobenzene	30.000	30.000	0.0	103	0.00
29	1,1-Dichloropropene	20.000	18.481	7.6	102	0.00
30 M	2-Butanone	100.000	96.897	3.1	100	0.00
31	2,2-Dichloropropane	20.000	19.241	3.8	103	0.00
32 M	1,1,1-Trichloroethane	20.000	18.586	7.1	101	0.00
33 M	Carbon Tetrachloride	20.000	18.054	9.7	97	0.00
34 M	Benzene	20.000	19.178	4.1	103	0.00
35	Methacrylonitrile	20.000	18.461	7.7	99	0.00
36 M	1,2-Dichloroethane	20.000	19.292	3.5	104	0.00
37 M	Trichloroethene	20.000	18.910	5.4	104	0.00
38	Methylcyclohexane	20.000	18.256	8.7	100	0.00
39 M	1,2-Dichloropropane	20.000	18.805	6.0	102	0.00
40	Dibromomethane	20.000	18.423	7.9	100	0.00
41 M	Bromodichloromethane	20.000	18.712	6.4	101	0.00
42 M	Vinyl Acetate	100.000	95.118	4.9	103	0.00
43	Ethyl Acetate	20.000	18.204	9.0	97	0.00
44	Isopropyl Acetate	20.000	18.414	7.9	101	0.00
45 T	1,4-Dioxane	400.000	374.339	6.4	103	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX111219\
 Data File : VX013374.D
 Acq On : 12 Nov 2019 09:58
 Operator : JC/SP
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 MSVOA_X
 LabSampleId :
 VSTDCCC020

Quant Time: Nov 12 12:36:20 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X111119W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Nov 12 01:16:13 2019
 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	18.523	7.4	103	0.00
47	n-amyl Acetate	20.000	18.525	7.4	104	0.00
48 M	t-1,3-Dichloropropene	20.000	18.687	6.6	106	0.00
49 T	cis-1,3-Dichloropropene	20.000	19.084	4.6	105	0.00
50 M	1,1,2-Trichloroethane	20.000	18.354	8.2	98	0.00
51	Ethyl methacrylate	20.000	18.480	7.6	104	0.00
52	1,3-Dichloropropane	20.000	18.927	5.4	101	0.00
53 M	Dibromochloromethane	20.000	18.721	6.4	104	0.00
54 M	1,2-Dibromoethane	20.000	18.844	5.8	101	0.00
55 M	2-Chloroethyl vinyl ether	100.000	92.771	7.2	99	0.00
56 M	Bromoform	20.000	18.039	9.8	104	0.00
57 I	Chlorobenzene-d5	30.000	30.000	0.0	104	0.00
58 M	4-Methyl-2-Pentanone	100.000	92.338	7.7	99	0.00
59 M	2-Hexanone	100.000	95.739	4.3	102	0.00
60 S	4-Bromofluorobenzene	30.000	30.247	-0.8	107	0.00
61 M	Tetrachloroethene	20.000	19.183	4.1	106	0.00
62 M	Toluene	20.000	19.033	4.8	102	0.00
63 S	Toluene-d8	30.000	30.516	-1.7	105	0.00
64 M	Chlorobenzene	20.000	19.131	4.3	106	0.00
65	1,1,1,2-Tetrachloroethane	20.000	18.595	7.0	103	0.00
66 M	Ethyl Benzene	20.000	18.911	5.4	103	0.00
67 M	m/p-Xylenes	40.000	37.854	5.4	104	0.00
68 M	o-Xylene	20.000	18.991	5.0	105	0.00
69 M	Styrene	20.000	18.823	5.9	104	0.00
70	Isopropylbenzene	20.000	19.034	4.8	104	0.00
71 M	1,1,2,2-Tetrachloroethane	20.000	18.661	6.7	102	0.00
72	1,2,3-Trichloropropane	20.000	20.861	-4.3	123	0.00
73	Bromobenzene	20.000	18.989	5.1	106	0.00
74	n-propylbenzene	20.000	19.024	4.9	105	0.00
75	2-Chlorotoluene	20.000	18.966	5.2	104	0.00
76	1,3,5-Trimethylbenzene	20.000	18.930	5.4	104	0.00
77	t-1,4-Dichloro-2-butene	20.000	17.704	11.5	103	0.00
78	4-Chlorotoluene	20.000	18.606	7.0	104	0.00
79	tert-butylbenzene	20.000	18.393	8.0	101	0.00
80	1,2,4-Trimethylbenzene	20.000	19.021	4.9	105	0.00
81	sec-Butylbenzene	20.000	18.882	5.6	105	0.00
82	p-Isopropyltoluene	20.000	18.994	5.0	106	0.00
83 M	1,3-Dichlorobenzene	20.000	19.137	4.3	109	0.00
84 M	1,4-Dichlorobenzene	20.000	18.925	5.4	108	0.00
85	n-Butylbenzene	20.000	18.557	7.2	106	0.00
86 T	Hexachloroethane	20.000	18.162	9.2	107	0.00
87 M	1,2-Dichlorobenzene	20.000	19.108	4.5	107	0.00
88	1,2-Dibromo-3-Chloropropane	20.000	17.606	12.0	102	0.00
89	1,2,4-Trichlorobenzene	20.000	18.629	6.9	107	0.00
90	Hexachlorobutadiene	20.000	18.184	9.1	104	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA X\Data\VX1111219\
Data File : VX013374.D
Acq On : 12 Nov 2019 09:58
Operator : JC/SP
Sample : VSTDCCC020
Misc : 5.0mL/MSVOA X/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument :
MSVOA_X
LabSampleId :
VSTDCCC020

Quant Time: Nov 12 12:36:20 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA X\METHOD\624X111119W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
QLast Update : Tue Nov 12 01:16:13 2019
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)
91 M	Naphthalene	20.000	18.609	7.0	105	0.00
92	1,2,3-Trichlorobenzene	20.000	18.502	7.5	105	0.00

(#) = Out of Range

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