

Data Path : Z:\voasrv\HPCHEM1\MSVOA N\Data\VN100620\
 Data File : VN063893.D
 Acq On : 6 Oct 2020 14:31
 Operator : JC/MD
 Sample : VSTDICV020
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 ICVVN100620

Quant Time: Oct 06 16:49:49 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100620W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Oct 06 16:42:38 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	0.00
2 M	Dichlorodifluoromethane	20.000	20.846	-4.2	108	0.00
3 M	Chloromethane	20.000	19.199	4.0	104	0.00
4 M	Vinyl Chloride	20.000	19.978	0.1	108	0.00
5 M	Bromomethane	20.000	19.322	3.4	107	0.02
6 M	Chloroethane	20.000	18.825	5.9	103	0.01
7 M	Trichlorofluoromethane	20.000	19.768	1.2	109	0.01
8 T	Diethyl Ether	20.000	19.032	4.8	108	0.00
9	1,1,2-Trichlorotrifluoroeth	20.000	19.992	0.0	109	0.00
10 M	1,1-Dichloroethene	20.000	19.377	3.1	108	0.00
11	Methyl Iodide	20.000	16.421	17.9	92	0.00
12	Methyl Acetate	20.000	17.511	12.4	97	0.00
13 M	Acrolein	100.000	92.814	7.2	128	0.00
14 M	Acrylonitrile	100.000	86.518	13.5	100	0.00
15 M	Acetone	100.000	93.824	6.2	105	0.00
16 M	Carbon Disulfide	20.000	19.493	2.5	110	0.00
17	Allyl chloride	20.000	19.176	4.1	101	0.00
18 M	Methylene Chloride	20.000	18.835	5.8	103	0.00
19 M	trans-1,2-Dichloroethene	20.000	19.409	3.0	113	0.00
20 T	Diisopropyl ether	20.000	19.243	3.8	109	0.00
21 M	1,1-Dichloroethane	20.000	19.228	3.9	106	0.00
22 M	cis-1,2-Dichloroethene	20.000	20.121	-0.6	117	0.00
23 M	tert-Butyl Alcohol	100.000	83.773	16.2	94	-0.02
24 M	Methyl tert-Butyl Ether	20.000	19.244	3.8	107	0.00
25 M	Chloroform	20.000	18.976	5.1	105	0.00
26	Cyclohexane	20.000	19.643	1.8	114	0.00
27 s	1,2-Dichloroethane-d4	30.000	30.927	-3.1	121	0.00
28 I	1,4-Difluorobenzene	30.000	30.000	0.0	124	0.00
29	1,1-Dichloropropene	20.000	18.242	8.8	106	0.00
30 M	2-Butanone	100.000	83.908	16.1	100	0.00
31	2,2-Dichloropropane	20.000	18.674	6.6	108	0.00
32 M	1,1,1-Trichloroethane	20.000	18.444	7.8	106	0.00
33 M	Carbon Tetrachloride	20.000	18.161	9.2	106	0.00
34 M	Benzene	20.000	18.411	7.9	107	0.00
35	Methacrylonitrile	20.000	14.877	25.6	94	0.00
36 M	1,2-Dichloroethane	20.000	17.582	12.1	103	0.00
37 M	Trichloroethene	20.000	18.265	8.7	110	0.00
38	Methylcyclohexane	20.000	18.989	5.1	114	0.00
39 M	1,2-Dichloropropane	20.000	18.385	8.1	105	0.00
40	Dibromomethane	20.000	17.553	12.2	101	0.00
41 M	Bromodichloromethane	20.000	18.100	9.5	107	0.00
42 M	Vinyl Acetate	100.000	88.416	11.6	104	0.00
43	Ethyl Acetate	20.000	16.521	17.4	101	0.00
44	Isopropyl Acetate	20.000	17.044	14.8	100	0.00
45 T	1,4-Dioxane	400.000	285.178	28.7	86	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA N\Data\VN100620\
 Data File : VN063893.D
 Acq On : 6 Oct 2020 14:31
 Operator : JC/MD
 Sample : VSTDICV020
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleID :
 ICVVN100620

Quant Time: Oct 06 16:49:49 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100620W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Tue Oct 06 16:42:38 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	16.657	16.7	97	0.00
47	n-amyl Acetate	20.000	15.937	20.3	99	0.00
48 M	t-1,3-Dichloropropene	20.000	17.778	11.1	105	0.00
49 T	cis-1,3-Dichloropropene	20.000	18.340	8.3	111	0.00
50 M	1,1,2-Trichloroethane	20.000	18.000	10.0	107	0.00
51	Ethyl methacrylate	20.000	17.165	14.2	104	0.00
52	1,3-Dichloropropane	20.000	17.826	10.9	107	0.00
53 M	Dibromochloromethane	20.000	17.721	11.4	104	0.00
54 M	1,2-Dibromoethane	20.000	17.833	10.8	105	0.00
55 M	2-Chloroethyl vinyl ether	100.000	58.018	42.0#	74	0.00
56 M	Bromoform	20.000	16.685	16.6	103	0.00
57 I	Chlorobenzene-d5	30.000	30.000	0.0	125	0.00
58 M	4-Methyl-2-Pentanone	100.000	83.221	16.8	97	0.00
59 M	2-Hexanone	100.000	81.737	18.3	97	0.00
60 S	4-Bromofluorobenzene	30.000	28.921	3.6	125	0.00
61 M	Tetrachloroethene	20.000	19.108	4.5	109	0.00
62 M	Toluene	20.000	18.086	9.6	108	0.00
63 S	Toluene-d8	30.000	29.955	0.2	124	0.00
64 M	Chlorobenzene	20.000	18.417	7.9	110	0.00
65	1,1,1,2-Tetrachloroethane	20.000	17.665	11.7	105	0.00
66 M	Ethyl Benzene	20.000	18.446	7.8	111	0.00
67 M	m/p-Xylenes	40.000	37.337	6.7	111	0.00
68 M	o-Xylene	20.000	17.902	10.5	109	0.00
69 M	Styrene	20.000	17.867	10.7	109	0.00
70	Isopropylbenzene	20.000	18.395	8.0	112	0.00
71 M	1,1,2,2-Tetrachloroethane	20.000	16.801	16.0	101	0.00
72	1,2,3-Trichloropropane	20.000	15.821	20.9	92	0.00
73	Bromobenzene	20.000	18.207	9.0	113	0.00
74	n-propylbenzene	20.000	18.459	7.7	112	0.00
75	2-Chlorotoluene	20.000	18.882	5.6	114	0.00
76	1,3,5-Trimethylbenzene	20.000	18.569	7.2	113	0.00
77	t-1,4-Dichloro-2-butene	20.000	15.885	20.6	100	0.00
78	4-Chlorotoluene	20.000	18.495	7.5	112	0.00
79	tert-butylbenzene	20.000	18.249	8.8	113	0.00
80	1,2,4-Trimethylbenzene	20.000	18.110	9.5	110	0.00
81	sec-Butylbenzene	20.000	18.232	8.8	111	0.00
82	p-Isopropyltoluene	20.000	18.091	9.5	113	0.00
83 M	1,3-Dichlorobenzene	20.000	18.179	9.1	112	0.00
84 M	1,4-Dichlorobenzene	20.000	18.120	9.4	113	0.00
85	n-Butylbenzene	20.000	18.451	7.7	116	0.00
86 T	Hexachloroethane	20.000	17.464	12.7	109	0.00
87 M	1,2-Dichlorobenzene	20.000	17.823	10.9	108	0.00
88	1,2-Dibromo-3-Chloropropane	20.000	15.743	21.3	95	0.00
89	1,2,4-Trichlorobenzene	20.000	18.073	9.6	114	0.00
90	Hexachlorobutadiene	20.000	19.345	3.3	115	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA N\Data\VN100620\
Data File : VN063893.D
Acq On : 6 Oct 2020 14:31
Operator : JC/MD
Sample : VSTDICV020
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
ICVVN100620

Quant Time: Oct 06 16:49:49 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA N\METHODS\624N100620W.M
Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
QLast Update : Tue Oct 06 16:42:38 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)
91 M	Naphthalene	20.000	17.163	14.2	105	0.00
92	1,2,3-Trichlorobenzene	20.000	18.192	9.0	112	0.00

(#) = Out of Range

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