

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061122\
 Data File : VN072794.D
 Acq On : 11 Jun 2022 17:43
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Jun 13 01:20:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N060622W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Mon Jun 06 14:10:15 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	Bromochloromethane	30.000	30.000	0.0	77	0.00
2 M	Dichlorodifluoromethane	20.000	16.356	18.2	63	0.00
3 M	Chloromethane	20.000	15.570	22.1	61	0.00
4 M	Vinyl Chloride	20.000	19.698	1.5	76	0.00
5 M	Bromomethane	20.000	14.190	29.1#	56	0.00
6 M	Chloroethane	20.000	19.963	0.2	72	0.00
7 M	Trichlorofluoromethane	20.000	17.276	13.6	66	0.00
8 T	Diethyl Ether	20.000	18.205	9.0	70	0.00
9	1,1,2-Trichlorotrifluoroeth	20.000	17.317	13.4	66	0.00
10 M	1,1-Dichloroethene	20.000	17.210	13.9	68	0.00
11	Methyl Iodide	20.000	10.174	49.1#	38	0.00
12	Methyl Acetate	20.000	20.237	-1.2	80	0.00
13 M	Acrolein	100.000	83.287	16.7	69	0.00
14 M	Acrylonitrile	100.000	98.346	1.7	75	0.00
15 M	Acetone	100.000	94.697	5.3	72	0.00
16 M	Carbon Disulfide	20.000	13.740	31.3#	56	0.00
17	Allyl chloride	20.000	17.501	12.5	68	0.00
18 M	Methylene Chloride	20.000	18.122	9.4	69	0.00
19 M	trans-1,2-Dichloroethene	20.000	16.897	15.5	65	0.00
20 T	Diisopropyl ether	20.000	20.386	-1.9	76	0.00
21 M	1,1-Dichloroethane	20.000	18.018	9.9	70	0.00
22 M	cis-1,2-Dichloroethene	20.000	18.414	7.9	70	0.00
23 M	tert-Butyl Alcohol	100.000	86.708	13.3	65	0.01
24 M	Methyl tert-Butyl Ether	20.000	18.354	8.2	70	0.01
25 M	Chloroform	20.000	18.871	5.6	72	0.00
26	Cyclohexane	20.000	17.411	12.9	67	0.00
27 s	1,2-Dichloroethane-d4	30.000	32.099	-7.0	80	0.00
28 I	1,4-Difluorobenzene	30.000	30.000	0.0	75	0.00
29	1,1-Dichloropropene	20.000	17.669	11.7	69	0.00
30 M	2-Butanone	100.000	101.030	-1.0	78	0.00
31	2,2-Dichloropropane	20.000	13.763	31.2#	53	0.00
32 M	1,1,1-Trichloroethane	20.000	17.191	14.0	66	0.00
33 M	Carbon Tetrachloride	20.000	16.409	18.0	63	0.00
34 M	Benzene	20.000	18.469	7.7	71	0.00
35	Methacrylonitrile	20.000	18.312	8.4	74	0.00
36 M	1,2-Dichloroethane	20.000	18.637	6.8	70	0.00
37 M	Trichloroethene	20.000	17.181	14.1	67	0.00
38	Methylcyclohexane	20.000	16.736	16.3	66	0.00
39 M	1,2-Dichloropropane	20.000	19.301	3.5	74	0.00
40	Dibromomethane	20.000	18.864	5.7	73	0.00
41 M	Bromodichloromethane	20.000	17.858	10.7	68	0.00
42 M	Vinyl Acetate	100.000	96.033	4.0	73	0.00
43	Ethyl Acetate	20.000	20.487	-2.4	77	0.00
44	Isopropyl Acetate	20.000	19.182	4.1	74	0.00
45 T	1,4-Dioxane	400.000	396.848	0.8	74	0.00
46	Methyl methacrylate	20.000	18.742	6.3	74	0.00
47	n-amyl Acetate	20.000	17.187	14.1	71	0.00
48 M	t-1,3-Dichloropropene	20.000	15.948	20.3	63	0.00

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061122\
 Data File : VN072794.D
 Acq On : 11 Jun 2022 17:43
 Operator : JC\MD
 Sample : VSTDCCC020
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_N
 LabSampleId :
 VSTDCCC020

Quant Time: Jun 13 01:20:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\624N060622W.M
 Quant Title : METHOD 624 VOLATILE ORGANIC ANALYSIS
 QLast Update : Mon Jun 06 14:10:15 2022
 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	Amount	Calc.	%Dev	Area%	Dev(min)
49 T	cis-1,3-Dichloropropene	20.000	16.680	16.6	65	0.00
50 M	1,1,2-Trichloroethane	20.000	18.713	6.4	71	0.00
51	Ethyl methacrylate	20.000	18.217	8.9	71	0.00
52	1,3-Dichloropropane	20.000	19.402	3.0	73	0.00
53 M	Dibromochloromethane	20.000	17.188	14.1	67	0.00
54 M	1,2-Dibromoethane	20.000	18.405	8.0	71	0.00
55 M	2-Chloroethyl vinyl ether	100.000	113.030	-13.0	94	0.00
56 M	Bromoform	20.000	15.687	21.6	63	0.00
57 I	Chlorobenzene-d5	30.000	30.000	0.0	74	0.00
58 M	4-Methyl-2-Pentanone	100.000	107.828	-7.8	80	0.00
59 M	2-Hexanone	100.000	103.366	-3.4	78	0.00
60 S	4-Bromofluorobenzene	30.000	27.595	8.0	70	0.00
61 M	Tetrachloroethene	20.000	18.887	5.6	70	0.00
62 M	Toluene	20.000	19.036	4.8	71	0.00
63 S	Toluene-d8	30.000	32.114	-7.0	77	0.00
64 M	Chlorobenzene	20.000	18.334	8.3	69	0.00
65	1,1,1,2-Tetrachloroethane	20.000	18.560	7.2	68	0.00
66 M	Ethyl Benzene	20.000	18.608	7.0	70	0.00
67 M	m/p-Xylenes	40.000	35.530	11.2	68	0.00
68 M	o-Xylene	20.000	17.989	10.1	67	0.00
69 M	Styrene	20.000	17.875	10.6	70	0.00
70	Isopropylbenzene	20.000	18.213	8.9	69	0.00
71 M	1,1,2,2-Tetrachloroethane	20.000	19.830	0.9	74	0.00
72	1,2,3-Trichloropropane	20.000	21.827	-9.1	74	0.00
73	Bromobenzene	20.000	17.188	14.1	66	0.00
74	n-propylbenzene	20.000	17.438	12.8	68	0.00
75	2-Chlorotoluene	20.000	18.107	9.5	70	0.00
76	1,3,5-Trimethylbenzene	20.000	18.117	9.4	70	0.00
77	t-1,4-Dichloro-2-butene	20.000	13.547	32.3#	54	0.00
78	4-Chlorotoluene	20.000	16.938	15.3	67	0.00
79	tert-butylbenzene	20.000	17.793	11.0	68	0.00
80	1,2,4-Trimethylbenzene	20.000	17.992	10.0	71	0.00
81	sec-Butylbenzene	20.000	17.992	10.0	71	0.00
82	p-Isopropyltoluene	20.000	17.197	14.0	70	0.00
83 M	1,3-Dichlorobenzene	20.000	17.137	14.3	70	0.00
84 M	1,4-Dichlorobenzene	20.000	17.118	14.4	69	0.00
85	n-Butylbenzene	20.000	16.125	19.4	67	0.00
86 T	Hexachloroethane	20.000	15.695	21.5	61	0.00
87 M	1,2-Dichlorobenzene	20.000	17.613	11.9	70	0.00
88	1,2-Dibromo-3-Chloropropane	20.000	16.878	15.6	65	0.00
89	1,2,4-Trichlorobenzene	20.000	15.409	23.0	64	0.00
90	Hexachlorobutadiene	20.000	16.465	17.7	63	0.00
91 M	Naphthalene	20.000	16.564	17.2	69	0.00
92	1,2,3-Trichlorobenzene	20.000	16.245	18.8	68	0.00

(#) = Out of Range

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