Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU073120\
Data File : VU039733.D
Acq On : 31 Jul 2020 14:55
Operator : SY/MD
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Aug 01 01:40:21 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U072920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Fri Jul 31 07:12:27 2020
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 30% Max. Rel. Area : 200%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	i	Fluorobenzene	1.000	1.000	0.0	114	0.00
2	T	Dichlorodifluoromethane	10.000	8.406	15.9	96	0.00
3	t	Chloromethane	10.000	8.410	15.9	101	0.00
4	Rt	Vinyl Chloride	10.000	8.745	12.6	104	0.00
5	T	Bromomethane	10.000	9.260	7.4	121	0.00
6	T	Chloroethane	10.000	11.393	-13.9	139	0.00
7	T	Trichlorofluoromethane	10.000	8.811	11.9	104	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	8.602	14.0	102	0.00
9	Rt	1,1-Dichloroethene	10.000	8.625	13.8	101	0.00
10	t	Iodomethane	10.000	11.516	-15.2	113	0.00
11	t	Allyl Chloride	10.000	8.375	16.3	100	0.00
12	t	Acrylonitrile	20.000	18.202	9.0	111	0.00
13	T	Acetone	51.000	40.423	20.7	98	0.01
14	T	Carbon Disulfide	10.000	8.508	14.9	100	0.00
15	RT	Methylene Chloride	10.000	7.687	23.1	103	0.00
16	RT	trans-1,2-Dichloroethene	10.000	8.702	13.0	104	0.00
17	t	1,1-Dichloroethane	10.000	8.559	14.4	101	0.00
18	T	2-Butanone	50.000	40.509	19.0	99	0.00
19		Cyclohexane	10.000	8.141	18.6	99	0.00
20		Methylcyclohexane	10.000	9.636	3.6	105	0.00
21	T	2,2-Dichloropropane	10.000	7.857	21.4	96	0.00
22	RT	cis-1,2-Dichloroethene	10.000	8.551	14.5	102	0.00
23	t	Diethyl Ether	10.000	8.666	13.3	104	0.00
24	t	tert-Butyl Alcohol	100.000	85.921	14.1	103	0.04
25	t	Methyl tert-Butyl Ether	10.000	8.684	13.2	101	0.00
26	t	Bromochloromethane	10.000	8.785	12.1	106	0.00
27	t	Chloroform	10.000	8.389	16.1	101	0.00
28	RT	1,1,1-Trichloroethane	10.000	8.461	15.4	101	0.00
29	T	1,1-Dichloropropene	10.000	8.350	16.5	100	0.00
30	RT	Carbon Tetrachloride	10.000	8.604	14.0	102	0.00
31	t	Isopropyl Ether	10.000	8.248	17.5	98	0.00
32		Ethyl-t-butyl ether	10.000	8.359	16.4	99	0.00
33		Tert-Amyl methyl ether	10.000	9.814	1.9	114	0.00
34	t	Propionitrile	50.000	45.234	9.5	108	0.00
35	RT	Benzene	10.000	9.366	6.3	112	0.00
36	RT	1,2-Dichloroethane	10.000	9.585	4.1	115	0.00
37	RT	Trichloroethene	10.000	8.880	11.2	104	0.00
38	Rt	1,2-Dichloropropane	10.000	9.422	5.8	107	0.00
39	t	Methacrylonitrile	10.000	7.763	22.4	103	0.00
40	t	Methyl acrylate	10.000	8.750	12.5	105	0.00
41	t	Tetrahydrofuran	20.000	16.414	17.9	108	0.00
42	t	1-Chlorobutane	10.000	8.301	17.0	99	0.00
43	T	Dibromomethane	10.000	9.171	8.3	110	0.00
44	T	Bromodichloromethane	10.000	8.916	10.8	100	0.00
45	T	4-Methyl-2-Pentanone	50.000	47.161	5.7	105	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	19.457	2.7	122	0.00

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Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U072920DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
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Max. RRF Dev : 30% Max. Rel. Area : 200%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
47	t	Methyl methacrylate	20.000	19.688	1.6	107	0.00
48	t	Ethyl methacrylate	10.000	9.368	6.3	102	0.00
49	Rt	Toluene	10.000	8.982	10.2	100	0.00
50	T	t-1,3-Dichloropropene	10.000	9.205	7.9	97	0.00
51	T	cis-1,3-Dichloropropene	10.000	8.927	10.7	98	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.132	8.7	107	0.00
53	t	1,3-Dichloropropane	10.000	8.943	10.6	104	0.00
54	t	2-Hexanone	50.000	47.075	5.8	105	0.00
55	t	Dibromochloromethane	10.000	9.093	9.1	100	0.00
56	T	1,2-Dibromoethane	10.000	8.938	10.6	101	0.00
57	S	4-Bromofluorobenzene	1.000	0.886	11.4	97	0.00
58	RT	Tetrachloroethene	10.000	8.676	13.2	102	0.00
59	Rt	Chlorobenzene	10.000	8.772	12.3	98	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	8.768	12.3	99	0.00
61	t	Pentachloroethane	10.000	8.823	11.8	97	0.00
62	t	Hexachloroethane	10.000	9.150	8.5	98	0.00
63	Rt	Ethyl Benzene	10.000	9.221	7.8	98	0.00
64	RT	m/p-Xylenes	20.000	18.950	5.3	100	0.00
65	RT	o-Xylene	10.000	9.305	7.0	100	0.00
66	RT	Styrene	10.000	9.613	3.9	100	0.00
67	t	Bromoform	10.000	9.579	4.2	101	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.929	7.1	103	0.00
69	T	Isopropylbenzene	10.000	9.126	8.7	98	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.150	8.5	104	0.00
71	T	1,2,3-Trichloropropane	10.000	9.736	2.6	99	0.00
72	t	Bromobenzene	10.000	8.796	12.0	99	0.00
73	t	n-propylbenzene	10.000	9.138	8.6	97	0.00
74	t	2-Chlorotoluene	10.000	9.050	9.5	100	0.00
75	t	1,3,5-Trimethylbenzene	10.000	9.443	5.6	98	0.00
76	t	4-Chlorotoluene	10.000	9.325	6.8	100	0.00
77	t	tert-Butylbenzene	10.000	9.265	7.3	98	0.00
78	t	1,2,4-Trimethylbenzene	10.000	9.814	1.9	101	0.00
79	t	sec-Butylbenzene	10.000	9.266	7.3	99	0.00
80		Nitrobenzene	50.000	48.822	2.4	110	0.00
81	t	p-Isopropyltoluene	10.000	9.448	5.5	99	0.00
82	t	1,3-Dichlorobenzene	10.000	8.775	12.2	99	0.00
83	Rt	1,4-Dichlorobenzene	10.000	8.707	12.9	99	0.00
84	t	n-Butylbenzene	10.000	9.204	8.0	98	0.00
85	Rt	1,2-Dichlorobenzene	10.000	8.766	12.3	99	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	9.713	2.9	104	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	8.556	14.4	93	0.00
88	t	Hexachlorobutadiene	10.000	8.582	14.2	98	0.00
89	t	Naphthalene	10.000	9.131	8.7	95	0.00
90	t	1,2,3-Trichlorobenzene	10.000	8.796	12.0	97	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0