Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU012220\
Data File : VU036530.D
Acq On : 22 Jan 2020 17:38
Operator : JC/MD
Sample : VSTDICV010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : ICVVU012220

Quant Time: Jan 23 09:22:41 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U012220DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Thu Jan 23 08:59:25 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	99	0.00
2	T	Dichlorodifluoromethane	10.000	7.740	22.6	77	0.00
3	t	Chloromethane	10.000	8.924	10.8	91	0.00
4	Rt	Vinyl Chloride	10.000	8.329	16.7	83	0.00
5	T	Bromomethane	10.000	8.762	12.4	98	0.00
6	T	Chloroethane	10.000	8.782	12.2	97	0.00
7	T	Trichlorofluoromethane	10.000	9.924	0.8	100	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	10.065	-0.6	101	0.00
9	Rt	1,1-Dichloroethene	10.000	10.005	-0.1	102	0.00
10	t	Iodomethane	10.000	8.894	11.1	93	0.00
11	t	Allyl Chloride	10.000	9.742	2.6	98	0.00
12	t	Acrylonitrile	20.000	46.936	-134.7#	240	0.00
13	T	Acetone	51.000	76.174	-49.4#	152	0.00
14	T	Carbon Disulfide	10.000	9.432	5.7	95	0.00
15	RT	Methylene Chloride	10.000	8.445	15.5	99	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.945	0.5	100	0.00
17	t	1,1-Dichloroethane	10.000	9.681	3.2	97	0.00
18	T	2-Butanone	50.000	54.322	-8.6	111	0.00
19		Cyclohexane	10.000	9.704	3.0	97	0.00
20		Methylcyclohexane	10.000	10.008	-0.1	98	0.00
21	T	2,2-Dichloropropane	10.000	9.724	2.8	96	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.108	-1.1	101	0.00
23	t	Diethyl Ether	10.000	9.174	8.3	91	0.00
24	t	tert-Butyl Alcohol	100.000	49.460	50.5#	50	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.522	4.8	95	0.00
26	t	Bromochloromethane	10.000	10.661	-6.6	107	0.00
27	t	Chloroform	10.000	9.707	2.9	97	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.815	1.9	97	0.00
29	T	1,1-Dichloropropene	10.000	9.610	3.9	96	0.00
30	RT	Carbon Tetrachloride	10.000	9.837	1.6	97	0.00
31	t	Isopropyl Ether	10.000	9.916	0.8	99	0.00
32		Ethyl-t-butyl ether	10.000	0.000	100.0#	0	-4.56#
33		Tert-Amyl methyl ether	10.000	0.000	100.0#	0	-5.99#
34	t	Propionitrile	50.000	106.363	-112.7#	195	0.00
35	RT	Benzene	10.000	9.772	2.3	98	0.00
36	RT	1,2-Dichloroethane	10.000	9.837	1.6	97	0.00
37	RT	Trichloroethene	10.000	9.803	2.0	98	0.00
38	Rt	1,2-Dichloropropane	10.000	9.645	3.6	97	0.00
39	t	Methacrylonitrile	10.000	9.839	1.6	97	0.00
40	t	Methyl acrylate	10.000	11.309	-13.1	102	0.00
41	t	Tetrahydrofuran	20.000	46.925	-134.6#	237	0.00
42	t	1-Chlorobutane	10.000	9.658	3.4	96	0.00
43	T	Dibromomethane	10.000	9.906	0.9	98	0.00
44	T	Bromodichloromethane	10.000	9.916	0.8	98	0.00
45	T	4-Methyl-2-Pentanone	50.000	47.905	4.2	96	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	11.000	45.0#	45	0.00

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Quant Time: Jan 23 09:22:41 2020
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U012220DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Thu Jan 23 08:59:25 2020
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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)
47	t	Methyl methacrylate	20.000	9.858	50.7#	49	0.00
48	t	Ethyl methacrylate	10.000	10.261	-2.6	99	0.00
49	Rt	Toluene	10.000	9.783	2.2	99	0.00
50	T	t-1,3-Dichloropropene	10.000	10.432	-4.3	99	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.373	-3.7	100	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.921	0.8	98	0.00
53	t	1,3-Dichloropropane	10.000	9.860	1.4	97	0.00
54	t	2-Hexanone	50.000	53.795	-7.6	105	0.00
55	t	Dibromochloromethane	10.000	10.214	-2.1	100	0.00
56	T	1,2-Dibromoethane	10.000	10.117	-1.2	100	0.00
57	S	4-Bromofluorobenzene	1.000	0.976	2.4	98	0.00
58	RT	Tetrachloroethene	10.000	9.770	2.3	101	0.00
59	Rt	Chlorobenzene	10.000	9.663	3.4	97	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.063	-0.6	99	0.00
61	t	Pentachloroethane	10.000	10.552	-5.5	99	0.00
62	t	Hexachloroethane	10.000	10.188	-1.9	99	0.00
63	Rt	Ethyl Benzene	10.000	10.072	-0.7	100	0.00
64	RT	m/p-Xylenes	20.000	19.802	1.0	99	0.00
65	RT	o-Xylene	10.000	9.771	2.3	97	0.00
66	RT	Styrene	10.000	9.944	0.6	99	0.00
67	t	Bromoform	10.000	10.201	-2.0	97	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.985	1.5	102	0.00
69	T	Isopropylbenzene	10.000	9.918	0.8	99	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.813	1.9	97	0.00
71	T	1,2,3-Trichloropropane	10.000	9.858	1.4	104	0.00
72	t	Bromobenzene	10.000	10.123	-1.2	100	0.00
73	t	n-propylbenzene	10.000	9.955	0.4	96	0.00
74	t	2-Chlorotoluene	10.000	9.713	2.9	98	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.167	-1.7	102	0.00
76	t	4-Chlorotoluene	10.000	9.958	0.4	98	0.00
77	t	tert-Butylbenzene	10.000	9.738	2.6	98	0.00
78	t	1,2,4-Trimethylbenzene	10.000	9.902	1.0	100	0.00
79	t	sec-Butylbenzene	10.000	9.185	8.1	93	0.00
80		Nitrobenzene	50.000	0.000	100.0#	0	-13.23#
81	t	p-Isopropyltoluene	10.000	10.166	-1.7	102	0.00
82	t	1,3-Dichlorobenzene	10.000	9.702	3.0	97	0.00
83	Rt	1,4-Dichlorobenzene	10.000	9.696	3.0	98	0.00
84	t	n-Butylbenzene	10.000	10.467	-4.7	102	0.00
85	Rt	1,2-Dichlorobenzene	10.000	9.622	3.8	98	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	9.582	4.2	99	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.233	-12.3	102	0.00
88	t	Hexachlorobutadiene	10.000	10.119	-1.2	102	0.00
89	t	Naphthalene	10.000	11.712	-17.1	102	0.00
90	t	1,2,3-Trichlorobenzene	10.000	10.864	-8.6	102	0.00

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

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