Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU072219\
Data File : VU033366.D
Acq On : 22 Jul 2019 21:52
Operator : JC/SP
Sample : VSTDICV010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : ICVVU072219

Quant Time: Jul 23 03:51:49 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U072219DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Tue Jul 23 03:43:51 2019
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	98	0.00
2	T	Dichlorodifluoromethane	10.000	5.723	42.8#	59	0.00
3	t	Chloromethane	10.000	7.540	24.6	80	0.00
4	Rt	Vinyl Chloride	10.000	7.930	20.7	83	0.00
5	T	Bromomethane	10.000	8.299	17.0	90	0.00
6	T	Chloroethane	10.000	8.398	16.0	81	0.00
7	T	Trichlorofluoromethane	10.000	8.601	14.0	90	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.570	4.3	102	0.00
9	Rt	1,1-Dichloroethene	10.000	10.093	-0.9	106	0.00
10	t	Iodomethane	10.000	9.471	5.3	93	0.00
11	t	Allyl Chloride	10.000	9.645	3.6	101	0.00
12	t	Acrylonitrile	20.000	47.252	-136.3#	258	0.00
13	T	Acetone	51.000	43.682	14.3	98	0.00
14	T	Carbon Disulfide	10.000	8.551	14.5	91	0.00
15	RT	Methylene Chloride	10.000	8.579	14.2	100	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.272	7.3	100	0.00
17	t	1,1-Dichloroethane	10.000	9.693	3.1	104	0.00
18	T	2-Butanone	50.000	50.342	-0.7	97	0.00
19		Cyclohexane	10.000	10.909	-9.1	104	0.00
20		Methylcyclohexane	10.000	10.995	-9.9	101	0.00
21	T	2,2-Dichloropropane	10.000	8.722	12.8	97	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.004	-0.0	101	0.00
23	t	Diethyl Ether	10.000	9.228	7.7	94	0.00
24	t	tert-Butyl Alcohol	100.000	49.152	50.8#	50	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.689	3.1	99	0.00
26	t	Bromochloromethane	10.000	9.947	0.5	102	0.00
27	t	Chloroform	10.000	8.812	11.9	100	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.759	2.4	101	0.00
29	T	1,1-Dichloropropene	10.000	10.105	-1.1	99	0.00
30	RT	Carbon Tetrachloride	10.000	9.750	2.5	100	0.00
31	t	Isopropyl Ether	10.000	11.274	-12.7	113	0.00
32		Ethyl-t-butyl ether	10.000	0.000	100.0#	0	-4.05#
33		Tert-Amyl methyl ether	10.000	0.000	100.0#	0	-5.56#
34	t	Propionitrile	50.000	114.094	-128.2#	206	0.00
35	RT	Benzene	10.000	10.247	-2.5	102	0.00
36	RT	1,2-Dichloroethane	10.000	9.929	0.7	101	0.00
37	RT	Trichloroethene	10.000	10.310	-3.1	104	0.00
38	Rt	1,2-Dichloropropane	10.000	10.075	-0.7	102	0.00
39	t	Methacrylonitrile	10.000	10.628	-6.3	102	0.00
40	t	Methyl acrylate	10.000	10.636	-6.4	97	0.00
41	t	Tetrahydrofuran	20.000	58.312	-191.6#	265	0.00
42	t	1-Chlorobutane	10.000	10.177	-1.8	99	0.00
43	T	Dibromomethane	10.000	9.887	1.1	100	0.00
44	T	Bromodichloromethane	10.000	10.060	-0.6	103	0.00
45	T	4-Methyl-2-Pentanone	50.000	54.626	-9.3	99	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	11.023	44.9#	48	0.00

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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	10.959	45.2#	49	0.00
48	t	Ethyl methacrylate	10.000	12.434	-24.3	110	0.00
49	Rt	Toluene	10.000	11.268	-12.7	103	0.00
50	T	t-1,3-Dichloropropene	10.000	11.227	-12.3	106	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.873	-8.7	107	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.225	-2.2	103	0.00
53	t	1,3-Dichloropropane	10.000	10.358	-3.6	103	0.00
54	t	2-Hexanone	50.000	55.714	-11.4	100	0.00
55	t	Dibromochloromethane	10.000	10.400	-4.0	102	0.00
56	T	1,2-Dibromoethane	10.000	10.420	-4.2	103	0.00
57	S	4-Bromofluorobenzene	1.000	1.055	-5.5	97	0.00
58	RT	Tetrachloroethene	10.000	10.490	-4.9	104	0.00
59	Rt	Chlorobenzene	10.000	10.598	-6.0	102	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.253	-2.5	105	0.00
61	t	Pentachloroethane	10.000	9.922	0.8	103	0.00
62	t	Hexachloroethane	10.000	10.328	-3.3	103	0.00
63	Rt	Ethyl Benzene	10.000	11.788	-17.9	105	0.00
64	RT	m/p-Xylenes	20.000	23.536	-17.7	105	0.00
65	RT	o-Xylene	10.000	11.404	-14.0	102	0.00
66	RT	Styrene	10.000	12.009	-20.1	103	0.00
67	t	Bromoform	10.000	11.149	-11.5	107	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.047	-4.7	103	0.00
69	T	Isopropylbenzene	10.000	12.033	-20.3	107	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.092	-0.9	102	0.00
71	T	1,2,3-Trichloropropane	10.000	9.757	2.4	103	0.00
72	t	Bromobenzene	10.000	11.141	-11.4	106	0.00
73	t	n-propylbenzene	10.000	11.585	-15.9	101	0.00
74	t	2-Chlorotoluene	10.000	11.249	-12.5	102	0.00
75	t	1,3,5-Trimethylbenzene	10.000	12.103	-21.0	104	0.00
76	t	4-Chlorotoluene	10.000	11.441	-14.4	103	0.00
77	t	tert-Butylbenzene	10.000	11.269	-12.7	102	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.972	-19.7	103	0.00
79	t	sec-Butylbenzene	10.000	10.620	-6.2	94	0.00
80		Nitrobenzene	50.000	8.519	83.0#	22	0.00
81	t	p-Isopropyltoluene	10.000	11.997	-20.0	103	0.00
82	t	1,3-Dichlorobenzene	10.000	10.386	-3.9	101	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.762	-7.6	102	0.00
84	t	n-Butylbenzene	10.000	11.549	-15.5	102	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.500	-5.0	102	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	11.389	-13.9	110	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	12.696	-27.0	114	0.00
88	t	Hexachlorobutadiene	10.000	10.678	-6.8	109	0.00
89	t	Naphthalene	10.000	10.873	-8.7	113	0.00
90	t	1,2,3-Trichlorobenzene	10.000	12.115	-21.2	109	0.00

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

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