Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU012919\
Data File : VU029309.D
Acq On : 29 Jan 2019 19:23
Operator : JC/SP
Sample : VSTDICV010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 9 Sample Multiplier: 1

Quant Time: Jan 30 12:36:07 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U012919DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Wed Jan 30 09:13:14 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	99	0.00
2	T	Dichlorodifluoromethane	10.000	5.206	47.9#	52	0.00
3	t	Chloromethane	10.000	7.758	22.4	80	0.00
4	Rt	Vinyl Chloride	10.000	8.218	17.8	84	0.00
5	T	Bromomethane	10.000	8.374	16.3	97	0.00
6	T	Chloroethane	10.000	9.077	9.2	92	0.00
7	T	Trichlorofluoromethane	10.000	8.950	10.5	92	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.790	2.1	99	0.00
9	Rt	1,1-Dichloroethene	10.000	9.690	3.1	99	0.00
10	t	Iodomethane	10.000	9.068	9.3	89	0.00
11	t	Allyl Chloride	10.000	9.640	3.6	100	0.00
12	t	Acrylonitrile	20.000	46.685	-133.4#	242	0.00
13	T	Acetone	51.000	46.159	9.5	108	0.00
14	T	Carbon Disulfide	10.000	9.180	8.2	97	0.00
15	RT	Methylene Chloride	10.000	8.363	16.4	95	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.931	0.7	103	0.00
17	t	1,1-Dichloroethane	10.000	9.312	6.9	98	0.00
18	T	2-Butanone	50.000	52.817	-5.6	105	0.00
19		Cyclohexane	10.000	10.912	-9.1	106	0.00
20		Methylcyclohexane	10.000	11.275	-12.8	105	0.00
21	T	2,2-Dichloropropane	10.000	9.310	6.9	95	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.898	1.0	99	0.00
23	t	Diethyl Ether	10.000	9.672	3.3	100	0.00
24	t	tert-Butyl Alcohol	100.000	47.273	52.7#	48	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.942	0.6	103	0.00
26	t	Bromochloromethane	10.000	10.873	-8.7	106	0.00
27	t	Chloroform	10.000	9.735	2.7	97	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.771	2.3	99	0.00
29	T	1,1-Dichloropropene	10.000	10.390	-3.9	99	0.00
30	RT	Carbon Tetrachloride	10.000	9.848	1.5	99	0.00
31	t	Isopropyl Ether	10.000	10.900	-9.0	109	0.00
32		Ethyl-t-butyl ether	10.000	0.000	100.0#	0	-4.06#
33		Tert-Amyl methyl ether	10.000	0.000	100.0#	0	-5.58#
34	t	Propionitrile	50.000	111.065	-122.1#	208	0.00
35	RT	Benzene	10.000	10.025	-0.3	98	0.00
36	RT	1,2-Dichloroethane	10.000	9.764	2.4	98	0.00
37	RT	Trichloroethene	10.000	9.856	1.4	96	0.00
38	Rt	1,2-Dichloropropane	10.000	9.910	0.9	98	0.00
39	t	Methacrylonitrile	10.000	10.352	-3.5	99	0.00
40	t	Methyl acrylate	10.000	10.527	-5.3	99	0.00
41	t	Tetrahydrofuran	20.000	56.739	-183.7#	269	0.00
42	t	1-Chlorobutane	10.000	10.047	-0.5	99	0.00
43	T	Dibromomethane	10.000	9.913	0.9	98	0.00
44	T	Bromodichloromethane	10.000	9.799	2.0	97	0.00
45	T	4-Methyl-2-Pentanone	50.000	52.060	-4.1	97	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	10.014	49.9#	50	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	11.366	43.2#	51	0.00
48	t	Ethyl methacrylate	10.000	11.149	-11.5	100	0.00
49	Rt	Toluene	10.000	10.938	-9.4	103	0.00
50	T	t-1,3-Dichloropropene	10.000	11.119	-11.2	103	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.602	-6.0	100	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.625	-6.3	102	0.00
53	t	1,3-Dichloropropane	10.000	10.221	-2.2	100	0.00
54	t	2-Hexanone	50.000	52.721	-5.4	98	0.00
55	t	Dibromochloromethane	10.000	9.912	0.9	97	0.00
56	T	1,2-Dibromoethane	10.000	10.873	-8.7	105	0.00
57	S	4-Bromofluorobenzene	1.000	1.072	-7.2	101	0.00
58	RT	Tetrachloroethene	10.000	10.345	-3.5	101	0.00
59	Rt	Chlorobenzene	10.000	10.631	-6.3	102	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	9.921	0.8	99	0.00
61	t	Pentachloroethane	10.000	10.029	-0.3	100	0.00
62	t	Hexachloroethane	10.000	10.400	-4.0	99	0.00
63	Rt	Ethyl Benzene	10.000	11.347	-13.5	103	0.00
64	RT	m/p-Xylenes	20.000	23.375	-16.9	105	0.00
65	RT	o-Xylene	10.000	11.174	-11.7	101	0.00
66	RT	Styrene	10.000	11.023	-10.2	97	0.00
67	t	Bromoform	10.000	10.463	-4.6	100	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.096	-9.6	107	0.00
69	T	Isopropylbenzene	10.000	11.713	-17.1	107	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	9.711	2.9	96	0.00
71	T	1,2,3-Trichloropropane	10.000	9.643	3.6	90	0.00
72	t	Bromobenzene	10.000	11.046	-10.5	105	0.00
73	t	n-propylbenzene	10.000	11.408	-14.1	102	0.00
74	t	2-Chlorotoluene	10.000	11.309	-13.1	102	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.555	-15.5	104	0.00
76	t	4-Chlorotoluene	10.000	11.137	-11.4	102	0.00
77	t	tert-Butylbenzene	10.000	11.102	-11.0	100	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.630	-16.3	102	0.00
79	t	sec-Butylbenzene	10.000	10.850	-8.5	96	0.00
80		Nitrobenzene	50.000	13.859	72.3#	22	0.00
81	t	p-Isopropyltoluene	10.000	11.513	-15.1	101	0.00
82	t	1,3-Dichlorobenzene	10.000	10.313	-3.1	101	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.531	-5.3	102	0.00
84	t	n-Butylbenzene	10.000	11.701	-17.0	103	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.585	-5.9	99	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	11.290	-12.9	102	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.972	-19.7	107	0.00
88	t	Hexachlorobutadiene	10.000	11.045	-10.4	102	0.00
89	t	Naphthalene	10.000	12.784	-27.8	105	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.754	-17.5	106	0.00

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

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