Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU100318\
Data File : VU027363.D
Acq On : 03 Oct 2018 17:34
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Oct 03 18:13:33 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U100318DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Wed Oct 03 17:14:05 2018
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	100	0.00
2	T	Dichlorodifluoromethane	10.000	10.115	-1.2	100	0.00
3	t	Chloromethane	10.000	10.280	-2.8	102	0.00
4	Rt	Vinyl Chloride	10.000	10.168	-1.7	102	0.00
5	T	Bromomethane	10.000	12.335	-23.4	118	0.00
6	T	Chloroethane	10.000	10.291	-2.9	105	0.00
7	T	Trichlorofluoromethane	10.000	10.037	-0.4	101	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.990	0.1	101	0.00
9	Rt	1,1-Dichloroethene	10.000	10.211	-2.1	104	0.00
10	t	Iodomethane	10.000	10.841	-8.4	93	0.00
11	t	Allyl Chloride	10.000	9.853	1.5	102	0.00
12	t	Acrylonitrile	20.000	18.583	7.1	95	0.00
13	T	Acetone	51.000	47.891	6.1	107	0.00
14	T	Carbon Disulfide	10.000	9.944	0.6	101	0.00
15	RT	Methylene Chloride	10.000	9.189	8.1	102	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.519	4.8	96	0.00
17	t	1,1-Dichloroethane	10.000	9.623	3.8	96	0.00
18	T	2-Butanone	50.000	49.838	0.3	106	0.00
19		Cyclohexane	10.000	10.105	-1.1	101	0.00
20		Methylcyclohexane	10.000	10.179	-1.8	97	0.00
21	T	2,2-Dichloropropane	10.000	9.979	0.2	105	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.374	-3.7	107	0.00
23	t	Diethyl Ether	10.000	10.090	-0.9	101	0.00
24	t	tert-Butyl Alcohol	100.000	92.614	7.4	95	0.00
25	t	Methyl tert-Butyl Ether	10.000	9.581	4.2	93	0.00
26	t	Bromochloromethane	10.000	10.784	-7.8	108	0.00
27	t	Chloroform	10.000	10.507	-5.1	107	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.464	-4.6	104	0.00
29	T	1,1-Dichloropropene	10.000	10.386	-3.9	105	0.00
30	RT	Carbon Tetrachloride	10.000	10.802	-8.0	106	0.00
31	t	Isopropyl Ether	10.000	9.232	7.7	94	0.00
32		Ethyl-t-butyl ether	10.000	10.243	-2.4	101	0.00
33		Tert-Amyl methyl ether	10.000	10.408	-4.1	109	0.00
34	t	Propionitrile	50.000	51.787	-3.6	102	0.00
35	RT	Benzene	10.000	10.507	-5.1	105	0.00
36	RT	1,2-Dichloroethane	10.000	10.516	-5.2	103	0.00
37	RT	Trichloroethene	10.000	10.318	-3.2	104	0.00
38	Rt	1,2-Dichloropropane	10.000	10.097	-1.0	101	0.00
39	t	Methacrylonitrile	10.000	9.567	4.3	103	0.00
40	t	Methyl acrylate	10.000	9.998	0.0	104	0.00
41	t	Tetrahydrofuran	20.000	17.876	10.6	103	0.00
42	t	1-Chlorobutane	10.000	10.371	-3.7	101	0.00
43	T	Dibromomethane	10.000	10.486	-4.9	99	0.00
44	T	Bromodichloromethane	10.000	10.456	-4.6	103	0.00
45	T	4-Methyl-2-Pentanone	50.000	53.830	-7.7	100	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	22.614	-13.1	100	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	20.765	-3.8	100	0.00
48	t	Ethyl methacrylate	10.000	10.782	-7.8	97	0.00
49	Rt	Toluene	10.000	10.534	-5.3	98	0.00
50	T	t-1,3-Dichloropropene	10.000	10.598	-6.0	100	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.287	-2.9	99	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.608	-6.1	105	0.00
53	t	1,3-Dichloropropane	10.000	10.371	-3.7	100	0.00
54	t	2-Hexanone	50.000	54.659	-9.3	99	0.00
55	t	Dibromochloromethane	10.000	10.888	-8.9	104	0.00
56	T	1,2-Dibromoethane	10.000	10.409	-4.1	101	0.00
57	S	4-Bromofluorobenzene	1.000	0.905	9.5	86	0.00
58	RT	Tetrachloroethene	10.000	9.991	0.1	97	0.00
59	Rt	Chlorobenzene	10.000	10.319	-3.2	98	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.617	-6.2	103	0.00
61	t	Pentachloroethane	10.000	10.853	-8.5	106	0.00
62	t	Hexachloroethane	10.000	11.482	-14.8	107	0.00
63	Rt	Ethyl Benzene	10.000	11.045	-10.4	98	0.00
64	RT	m/p-Xylenes	20.000	23.359	-16.8	101	0.00
65	RT	o-Xylene	10.000	11.145	-11.4	97	0.00
66	RT	Styrene	10.000	11.762	-17.6	100	0.00
67	t	Bromoform	10.000	10.965	-9.6	103	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.049	-4.9	95	0.00
69	T	Isopropylbenzene	10.000	11.307	-13.1	97	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.491	-4.9	102	0.00
71	T	1,2,3-Trichloropropane	10.000	9.592	4.1	92	0.00
72	t	Bromobenzene	10.000	10.844	-8.4	102	0.00
73	t	n-propylbenzene	10.000	11.997	-20.0	100	0.00
74	t	2-Chlorotoluene	10.000	11.466	-14.7	100	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.826	-18.3	100	0.00
76	t	4-Chlorotoluene	10.000	11.627	-16.3	100	0.00
77	t	tert-Butylbenzene	10.000	11.417	-14.2	100	0.00
78	t	1,2,4-Trimethylbenzene	10.000	12.078	-20.8	101	0.00
79	t	sec-Butylbenzene	10.000	11.758	-17.6	99	0.00
80		Nitrobenzene	50.000	62.294	-24.6	113	0.00
81	t	p-Isopropyltoluene	10.000	12.220	-22.2	100	0.00
82	t	1,3-Dichlorobenzene	10.000	11.158	-11.6	102	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.249	-12.5	102	0.00
84	t	n-Butylbenzene	10.000	12.296	-23.0	100	0.00
85	Rt	1,2-Dichlorobenzene	10.000	11.168	-11.7	102	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	11.147	-11.5	103	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.415	-14.1	100	0.00
88	t	Hexachlorobutadiene	10.000	11.212	-12.1	105	0.00
89	t	Naphthalene	10.000	11.703	-17.0	97	0.00
90	t	1,2,3-Trichlorobenzene	10.000	11.653	-16.5	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