Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU121418\
Data File : VU028666.D
Acq On : 13 Dec 2018 09:08
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Dec 14 05:40:29 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U112618DW.M
Quant Title : METHOD 524.2 VOLATILES DRINKING WATER
QLast Update : Tue Nov 27 01:03:50 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	32	0.00
2	T	Dichlorodifluoromethane	10.000	10.550	-5.5	34	0.00
3	t	Chloromethane	10.000	9.517	4.8	32	0.00
4	Rt	Vinyl Chloride	10.000	10.827	-8.3	34	0.00
5	T	Bromomethane	10.000	10.524	-5.2	38	-0.02
6	T	Chloroethane	10.000	10.519	-5.2	35	0.00
7	T	Trichlorofluoromethane	10.000	10.802	-8.0	35	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	11.057	-10.6	35	0.00
9	Rt	1,1-Dichloroethene	10.000	10.603	-6.0	35	0.00
10	t	Iodomethane	10.000	17.476	-74.8#	51	0.00
11	t	Allyl Chloride	10.000	10.642	-6.4	34	0.00
12	t	Acrylonitrile	20.000	21.243	-6.2	36	0.00
13	T	Acetone	50.000	50.920	-1.8	36	-0.01
14	T	Carbon Disulfide	10.000	10.480	-4.8	34	0.00
15	RT	Methylene Chloride	10.000	10.843	-8.4	36	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.763	-7.6	35	0.00
17	t	1,1-Dichloroethane	10.000	10.743	-7.4	35	0.00
18	T	2-Butanone	50.000	55.923	-11.8	34	0.00
19		Cyclohexane	10.000	10.502	-5.0	33	0.00
20		Methylcyclohexane	10.000	9.889	1.1	30	0.00
21	T	2,2-Dichloropropane	10.000	10.890	-8.9	35	0.00
22	RT	cis-1,2-Dichloroethene	10.000	10.471	-4.7	32	0.00
23	t	Diethyl Ether	10.000	10.558	-5.6	34	0.00
24	t	tert-Butyl Alcohol	100.000	106.309	-6.3	33	-0.03
25	t	Methyl tert-Butyl Ether	10.000	10.758	-7.6	34	0.00
26	t	Bromochloromethane	10.000	10.275	-2.8	32	0.00
27	t	Chloroform	10.000	9.799	2.0	33	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.371	-3.7	33	0.00
29	T	1,1-Dichloropropene	10.000	10.237	-2.4	32	0.00
30	RT	Carbon Tetrachloride	10.000	10.334	-3.3	34	0.00
31	t	Isopropyl Ether	10.000	10.827	-8.3	34	0.00
32		Ethyl-t-butyl ether	10.000	10.727	-7.3	34	0.00
33		Tert-Amyl methyl ether	10.000	10.189	-1.9	32	0.00
34	t	Propionitrile	50.000	57.784	-15.6	35	-0.01
35	RT	Benzene	10.000	9.791	2.1	31	0.00
36	RT	1,2-Dichloroethane	10.000	10.287	-2.9	33	0.00
37	RT	Trichloroethene	10.000	10.052	-0.5	32	0.00
38	Rt	1,2-Dichloropropane	10.000	10.160	-1.6	32	0.00
39	t	Methacrylonitrile	10.000	10.875	-8.8	35	0.00
40	t	Methyl acrylate	10.000	10.466	-4.7	34	0.00
41	t	Tetrahydrofuran	20.000	21.834	-9.2	35	0.00
42	t	1-Chlorobutane	10.000	9.990	0.1	31	0.00
43	T	Dibromomethane	10.000	10.161	-1.6	32	0.00
44	T	Bromodichloromethane	10.000	9.937	0.6	32	0.00
45	T	4-Methyl-2-Pentanone	50.000	53.405	-6.8	33	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	21.236	-6.2	31	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	21.189	-5.9	33	0.00
48	t	Ethyl methacrylate	10.000	11.046	-10.5	33	0.00
49	Rt	Toluene	10.000	10.517	-5.2	32	0.00
50	T	t-1,3-Dichloropropene	10.000	10.575	-5.7	33	0.00
51	T	cis-1,3-Dichloropropene	10.000	10.488	-4.9	33	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.479	-4.8	33	0.00
53	t	1,3-Dichloropropane	10.000	10.436	-4.4	33	0.00
54	t	2-Hexanone	50.000	53.605	-7.2	33	0.00
55	t	Dibromochloromethane	10.000	10.845	-8.5	33	0.00
56	T	1,2-Dibromoethane	10.000	10.797	-8.0	35	0.00
57	S	4-Bromofluorobenzene	1.000	1.088	-8.8	33	0.00
58	RT	Tetrachloroethene	10.000	10.221	-2.2	32	0.00
59	Rt	Chlorobenzene	10.000	10.428	-4.3	33	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.926	-9.3	35	0.00
61	t	Pentachloroethane	10.000	11.367	-13.7	36	0.00
62	t	Hexachloroethane	10.000	12.024	-20.2	37	0.00
63	Rt	Ethyl Benzene	10.000	10.807	-8.1	33	0.00
64	RT	m/p-Xylenes	20.000	22.231	-11.2	34	0.00
65	RT	o-Xylene	10.000	11.206	-12.1	34	0.00
66	RT	Styrene	10.000	11.490	-14.9	35	0.00
67	t	Bromoform	10.000	11.250	-12.5	35	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.093	-9.3	35	0.00
69	T	Isopropylbenzene	10.000	11.060	-10.6	33	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	11.052	-10.5	34	0.00
71	T	1,2,3-Trichloropropane	10.000	10.007	-0.1	31	0.00
72	t	Bromobenzene	10.000	11.249	-12.5	34	0.00
73	t	n-propylbenzene	10.000	11.633	-16.3	34	0.00
74	t	2-Chlorotoluene	10.000	11.550	-15.5	36	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.518	-15.2	34	0.00
76	t	4-Chlorotoluene	10.000	11.710	-17.1	35	0.00
77	t	tert-Butylbenzene	10.000	11.572	-15.7	35	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.640	-16.4	35	0.00
79	t	sec-Butylbenzene	10.000	11.533	-15.3	34	0.00
80		Nitrobenzene	50.000	59.474	-18.9	35	0.00
81	t	p-Isopropyltoluene	10.000	11.559	-15.6	34	0.00
82	t	1,3-Dichlorobenzene	10.000	11.606	-16.1	35	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.655	-16.5	35	0.00
84	t	n-Butylbenzene	10.000	11.749	-17.5	34	0.00
85	Rt	1,2-Dichlorobenzene	10.000	11.529	-15.3	36	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.661	-6.6	33	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	11.311	-13.1	33	0.00
88	t	Hexachlorobutadiene	10.000	11.563	-15.6	35	0.00
89	t	Naphthalene	10.000	10.919	-9.2	33	0.00
90	t	1,2,3-Trichlorobenzene	10.000	10.998	-10.0	33	0.00

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

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