Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU121418\
Data File : VU028678.D
Acq On : 13 Dec 2018 15:07
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Dec 14 06:05:51 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	29	0.00
2	T	Dichlorodifluoromethane	10.000	10.499	-5.0	30	0.00
3	t	Chloromethane	10.000	9.356	6.4	28	0.00
4	Rt	Vinyl Chloride	10.000	10.569	-5.7	30	0.00
5	T	Bromomethane	10.000	10.243	-2.4	33	-0.02
6	T	Chloroethane	10.000	10.278	-2.8	30	0.00
7	T	Trichlorofluoromethane	10.000	10.763	-7.6	31	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	11.038	-10.4	31	0.00
9	Rt	1,1-Dichloroethene	10.000	10.524	-5.2	31	0.00
10	t	Iodomethane	10.000	17.601	-76.0#	46	0.00
11	t	Allyl Chloride	10.000	10.005	-0.1	29	0.00
12	t	Acrylonitrile	20.000	21.172	-5.9	32	0.00
13	T	Acetone	50.000	47.629	4.7	30	-0.01
14	T	Carbon Disulfide	10.000	10.052	-0.5	29	0.00
15	RT	Methylene Chloride	10.000	10.393	-3.9	31	0.00
16	RT	trans-1,2-Dichloroethene	10.000	10.364	-3.6	30	0.00
17	t	1,1-Dichloroethane	10.000	10.411	-4.1	30	0.00
18	T	2-Butanone	50.000	52.814	-5.6	29	0.00
19		Cyclohexane	10.000	10.065	-0.6	28	0.00
20		Methylcyclohexane	10.000	9.510	4.9	26	0.00
21	T	2,2-Dichloropropane	10.000	10.211	-2.1	29	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.967	0.3	28	0.00
23	t	Diethyl Ether	10.000	10.539	-5.4	30	0.00
24	t	tert-Butyl Alcohol	100.000	104.136	-4.1	29	-0.03
25	t	Methyl tert-Butyl Ether	10.000	10.468	-4.7	30	0.00
26	t	Bromochloromethane	10.000	10.056	-0.6	28	0.00
27	t	Chloroform	10.000	9.511	4.9	29	0.00
28	RT	1,1,1-Trichloroethane	10.000	10.363	-3.6	30	0.00
29	T	1,1-Dichloropropene	10.000	9.741	2.6	28	0.00
30	RT	Carbon Tetrachloride	10.000	9.970	0.3	29	0.00
31	t	Isopropyl Ether	10.000	10.373	-3.7	30	0.00
32		Ethyl-t-butyl ether	10.000	10.208	-2.1	29	0.00
33		Tert-Amyl methyl ether	10.000	10.009	-0.1	28	0.00
34	t	Propionitrile	50.000	56.826	-13.7	31	-0.01
35	RT	Benzene	10.000	9.623	3.8	27	0.00
36	RT	1,2-Dichloroethane	10.000	10.062	-0.6	29	0.00
37	RT	Trichloroethene	10.000	9.779	2.2	28	0.00
38	Rt	1,2-Dichloropropane	10.000	9.860	1.4	28	0.00
39	t	Methacrylonitrile	10.000	10.378	-3.8	30	0.00
40	t	Methyl acrylate	10.000	10.294	-2.9	30	0.00
41	t	Tetrahydrofuran	20.000	20.885	-4.4	30	0.00
42	t	1-Chlorobutane	10.000	9.663	3.4	27	0.00
43	T	Dibromomethane	10.000	9.906	0.9	28	0.00
44	T	Bromodichloromethane	10.000	9.662	3.4	28	0.00
45	T	4-Methyl-2-Pentanone	50.000	51.678	-3.4	29	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	19.892	0.5	26	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU121418\
Data File : VU028678.D
Acq On : 13 Dec 2018 15:07
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Dec 14 06:05:51 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)
47	t	Methyl methacrylate	20.000	21.166	-5.8	29	0.00
48	t	Ethyl methacrylate	10.000	10.196	-2.0	28	0.00
49	Rt	Toluene	10.000	10.019	-0.2	27	0.00
50	T	t-1,3-Dichloropropene	10.000	9.842	1.6	27	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.683	3.2	27	0.00
52	RT	1,1,2-Trichloroethane	10.000	10.446	-4.5	30	0.00
53	t	1,3-Dichloropropane	10.000	10.204	-2.0	29	0.00
54	t	2-Hexanone	50.000	50.374	-0.7	28	0.00
55	t	Dibromochloromethane	10.000	10.524	-5.2	29	0.00
56	T	1,2-Dibromoethane	10.000	10.409	-4.1	30	0.00
57	S	4-Bromofluorobenzene	1.000	1.040	-4.0	28	0.00
58	RT	Tetrachloroethene	10.000	10.420	-4.2	30	0.00
59	Rt	Chlorobenzene	10.000	10.288	-2.9	29	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	10.625	-6.3	30	0.00
61	t	Pentachloroethane	10.000	11.166	-11.7	32	0.00
62	t	Hexachloroethane	10.000	11.635	-16.3	33	0.00
63	Rt	Ethyl Benzene	10.000	10.508	-5.1	29	0.00
64	RT	m/p-Xylenes	20.000	21.905	-9.5	30	0.00
65	RT	o-Xylene	10.000	10.742	-7.4	29	0.00
66	RT	Styrene	10.000	11.011	-10.1	30	0.00
67	t	Bromoform	10.000	10.533	-5.3	29	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	1.136	-13.6	33	0.00
69	T	Isopropylbenzene	10.000	10.729	-7.3	29	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	11.021	-10.2	31	0.00
71	T	1,2,3-Trichloropropane	10.000	10.305	-3.0	29	0.00
72	t	Bromobenzene	10.000	11.254	-12.5	31	0.00
73	t	n-propylbenzene	10.000	11.344	-13.4	30	0.00
74	t	2-Chlorotoluene	10.000	11.355	-13.6	31	0.00
75	t	1,3,5-Trimethylbenzene	10.000	11.217	-12.2	30	0.00
76	t	4-Chlorotoluene	10.000	11.402	-14.0	30	0.00
77	t	tert-Butylbenzene	10.000	11.243	-12.4	30	0.00
78	t	1,2,4-Trimethylbenzene	10.000	11.296	-13.0	31	0.00
79	t	sec-Butylbenzene	10.000	11.146	-11.5	30	0.00
80		Nitrobenzene	50.000	53.769	-7.5	28	0.00
81	t	p-Isopropyltoluene	10.000	11.205	-12.1	30	0.00
82	t	1,3-Dichlorobenzene	10.000	11.357	-13.6	31	0.00
83	Rt	1,4-Dichlorobenzene	10.000	11.429	-14.3	31	0.00
84	t	n-Butylbenzene	10.000	11.256	-12.6	29	0.00
85	Rt	1,2-Dichlorobenzene	10.000	11.163	-11.6	31	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.224	-2.2	29	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	10.696	-7.0	28	0.00
88	t	Hexachlorobutadiene	10.000	11.555	-15.5	32	0.00
89	t	Naphthalene	10.000	10.803	-8.0	29	0.00
90	t	1,2,3-Trichlorobenzene	10.000	10.843	-8.4	29	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU121418\
Data File : VU028678.D
Acq On : 13 Dec 2018 15:07
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Dec 14 06:05:51 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)

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