Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU112618\
Data File : VU028340.D
Acq On : 26 Nov 2018 15:45
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Nov 27 01:52:36 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U0112618DW.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	102	0.00
2	T	Dichlorodifluoromethane	10.000	9.498	5.0	96	0.00
3	t	Chloromethane	10.000	9.133	8.7	98	0.00
4	Rt	Vinyl Chloride	10.000	9.723	2.8	97	0.00
5	T	Bromomethane	10.000	8.047	19.5	91	0.00
6	T	Chloroethane	10.000	9.582	4.2	100	0.00
7	T	Trichlorofluoromethane	10.000	9.793	2.1	99	0.00
8		1,1,2-Trichloro-1,2,2-trifl	10.000	9.972	0.3	100	0.00
9	Rt	1,1-Dichloroethene	10.000	9.564	4.4	100	0.00
10	t	Iodomethane	10.000	11.785	-17.9	109	0.00
11	t	Allyl Chloride	10.000	9.722	2.8	98	0.00
12	t	Acrylonitrile	20.000	20.034	-0.2	108	0.00
13	T	Acetone	50.000	43.857	12.3	99	0.00
14	T	Carbon Disulfide	10.000	9.574	4.3	98	0.00
15	RT	Methylene Chloride	10.000	9.578	4.2	101	0.00
16	RT	trans-1,2-Dichloroethene	10.000	9.572	4.3	99	0.00
17	t	1,1-Dichloroethane	10.000	9.821	1.8	100	0.00
18	T	2-Butanone	50.000	50.021	-0.0	97	0.00
19		Cyclohexane	10.000	10.116	-1.2	100	0.00
20		Methylcyclohexane	10.000	10.250	-2.5	100	0.00
21	T	2,2-Dichloropropane	10.000	9.480	5.2	97	0.00
22	RT	cis-1,2-Dichloroethene	10.000	9.778	2.2	96	0.00
23	t	Diethyl Ether	10.000	9.828	1.7	99	0.00
24	t	tert-Butyl Alcohol	100.000	94.602	5.4	93	0.02
25	t	Methyl tert-Butyl Ether	10.000	10.012	-0.1	100	0.00
26	t	Bromochloromethane	10.000	9.626	3.7	96	0.00
27	t	Chloroform	10.000	9.403	6.0	100	0.00
28	RT	1,1,1-Trichloroethane	10.000	9.954	0.5	102	0.00
29	T	1,1-Dichloropropene	10.000	9.965	0.4	99	0.00
30	RT	Carbon Tetrachloride	10.000	9.619	3.8	100	0.00
31	t	Isopropyl Ether	10.000	9.902	1.0	100	0.00
32		Ethyl-t-butyl ether	10.000	9.955	0.4	100	0.00
33		Tert-Amyl methyl ether	10.000	9.986	0.1	99	0.00
34	t	Propionitrile	50.000	53.047	-6.1	103	0.01
35	RT	Benzene	10.000	9.939	0.6	99	0.00
36	RT	1,2-Dichloroethane	10.000	9.822	1.8	100	0.00
37	RT	Trichloroethene	10.000	9.959	0.4	100	0.00
38	Rt	1,2-Dichloropropane	10.000	9.938	0.6	99	0.00
39	t	Methacrylonitrile	10.000	9.742	2.6	99	0.00
40	t	Methyl acrylate	10.000	9.445	5.5	96	0.00
41	t	Tetrahydrofuran	20.000	19.743	1.3	99	0.00
42	t	1-Chlorobutane	10.000	9.942	0.6	99	0.00
43	T	Dibromomethane	10.000	9.995	0.1	99	0.00
44	T	Bromodichloromethane	10.000	9.922	0.8	100	0.00
45	T	4-Methyl-2-Pentanone	50.000	51.089	-2.2	99	0.00
46	t	t-1,4-Dichloro-2-butene	20.000	20.946	-4.7	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU112618\
Data File : VU028340.D
Acq On : 26 Nov 2018 15:45
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTDCCC010

Quant Time: Nov 27 01:52:36 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U0112618DW.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	20.812	-4.1	101	0.00
48	t	Ethyl methacrylate	10.000	10.306	-3.1	99	0.00
49	Rt	Toluene	10.000	10.121	-1.2	98	0.00
50	T	t-1,3-Dichloropropene	10.000	10.011	-0.1	98	0.00
51	T	cis-1,3-Dichloropropene	10.000	9.976	0.2	99	0.00
52	RT	1,1,2-Trichloroethane	10.000	9.673	3.3	98	0.00
53	t	1,3-Dichloropropane	10.000	9.966	0.3	100	0.00
54	t	2-Hexanone	50.000	49.976	0.0	98	0.00
55	t	Dibromochloromethane	10.000	9.715	2.9	94	0.00
56	T	1,2-Dibromoethane	10.000	10.065	-0.6	102	0.00
57	S	4-Bromofluorobenzene	1.000	0.930	7.0	90	0.00
58	RT	Tetrachloroethene	10.000	9.741	2.6	98	0.00
59	Rt	Chlorobenzene	10.000	9.855	1.4	98	0.00
60	T	1,1,1,2-Tetrachloroethane	10.000	9.601	4.0	97	0.00
61	t	Pentachloroethane	10.000	9.591	4.1	96	0.00
62	t	Hexachloroethane	10.000	9.518	4.8	94	0.00
63	Rt	Ethyl Benzene	10.000	10.253	-2.5	99	0.00
64	RT	m/p-Xylenes	20.000	20.540	-2.7	99	0.00
65	RT	o-Xylene	10.000	10.232	-2.3	97	0.00
66	RT	Styrene	10.000	10.369	-3.7	99	0.00
67	t	Bromoform	10.000	9.728	2.7	95	0.00
68	S	1,2-Dichlorobenzene-d4	1.000	0.970	3.0	98	0.00
69	T	Isopropylbenzene	10.000	10.210	-2.1	98	0.00
70	T	1,1,2,2-Tetrachloroethane	10.000	10.117	-1.2	100	0.00
71	T	1,2,3-Trichloropropane	10.000	10.038	-0.4	100	0.00
72	t	Bromobenzene	10.000	10.193	-1.9	98	0.00
73	t	n-propylbenzene	10.000	10.539	-5.4	98	0.00
74	t	2-Chlorotoluene	10.000	10.216	-2.2	100	0.00
75	t	1,3,5-Trimethylbenzene	10.000	10.461	-4.6	98	0.00
76	t	4-Chlorotoluene	10.000	10.404	-4.0	97	0.00
77	t	tert-Butylbenzene	10.000	10.277	-2.8	98	0.00
78	t	1,2,4-Trimethylbenzene	10.000	10.502	-5.0	100	0.00
79	t	sec-Butylbenzene	10.000	10.571	-5.7	100	0.00
80		Nitrobenzene	50.000	53.717	-7.4	100	0.00
81	t	p-Isopropyltoluene	10.000	10.466	-4.7	97	0.00
82	t	1,3-Dichlorobenzene	10.000	10.283	-2.8	99	0.00
83	Rt	1,4-Dichlorobenzene	10.000	10.134	-1.3	97	0.00
84	t	n-Butylbenzene	10.000	10.849	-8.5	99	0.00
85	Rt	1,2-Dichlorobenzene	10.000	10.083	-0.8	98	0.00
86	t	1,2-Dibromo-3-Chloropropane	10.000	10.340	-3.4	102	0.00
87	Rt	1,2,4-Trichlorobenzene	10.000	10.551	-5.5	98	0.00
88	t	Hexachlorobutadiene	10.000	10.057	-0.6	97	0.00
89	t	Naphthalene	10.000	10.424	-4.2	98	0.00
90	t	1,2,3-Trichlorobenzene	10.000	10.153	-1.5	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU112618\
Data File : VU028340.D
Acq On : 26 Nov 2018 15:45
Operator : MD/SY
Sample : VSTDCCC010
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
LabSampleId :
VSTDCCC010

Quant Time: Nov 27 01:52:36 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\524U0112618DW.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