Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU010419\
Data File : VU029015.D
Acq On : 04 Jan 2019 09:37
Operator : MD/SY
Sample : VSTDCCC005
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTD00551

Quant Time: Jan 04 22:14:04 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR122818WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Thu Jan 03 03:47:56 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	1,4-Difluorobenzene	5.000	5.000	0.0	71	0.00
2	T	Dichlorodifluoromethane	5.000	5.013	-0.3	78	0.00
3	T	Chloromethane	5.000	5.098	-2.0	80	0.00
4	S	Vinyl Chloride-d3	5.000	6.066	-21.3	90	0.00
5	T	Vinyl chloride	5.000	5.130	-2.6	78	0.00
6	T	Bromomethane	5.000	4.865	2.7	77	0.00
7	S	Chloroethane-d5	5.000	5.736	-14.7	86	0.00
8	T	Chloroethane	5.000	5.046	-0.9	78	0.00
9	T	Trichlorofluoromethane	5.000	4.662	6.8	71	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	5.000	4.598	8.0	71	0.00
11	S	1,1-Dichloroethene-d2	5.000	5.282	-5.6	82	0.00
12	T	1,1-Dichloroethene	5.000	4.501	10.0	69	0.00
13	T	Acetone	50.000	47.945	4.1	74	0.00
14	T	Carbon disulfide	5.000	4.790	4.2	74	0.00
15	T	Methyl Acetate	5.000	4.963	0.7	81	0.00
16	T	Methylene chloride	5.000	3.670	26.6	69	0.00
17	T	Methyl tert-butyl Ether	5.000	4.791	4.2	72	0.00
18	T	trans-1,2-Dichloroethene	5.000	4.508	9.8	69	0.00
19	T	1,1-Dichloroethane	5.000	5.020	-0.4	76	0.00
20	S	2-Butanone-d5	50.000	57.736	-15.5	83	0.00
21	T	2-Butanone	50.000	56.767	-13.5	85	0.00
22	T	cis-1,2-Dichloroethene	5.000	4.926	1.5	75	0.00
23	T	Bromochloromethane	5.000	5.084	-1.7	77	0.00
24	S	Chloroform-d	5.000	5.596	-11.9	83	0.00
25	T	Chloroform	5.000	5.016	-0.3	78	0.00
26	S	1,2-Dichloroethane-d4	5.000	5.350	-7.0	81	0.00
27	T	1,2-Dichloroethane	5.000	5.144	-2.9	79	0.00
28	I	Chlorobenzene-d5	5.000	5.000	0.0	72	0.00
29	T	1,1,1-Trichloroethane	5.000	5.081	-1.6	77	0.00
30	T	Cyclohexane	5.000	4.973	0.5	76	0.00
31	T	Carbon tetrachloride	5.000	5.123	-2.5	78	0.00
32	S	Benzene-d6	5.000	5.528	-10.6	83	0.00
33	T	Benzene	5.000	5.052	-1.0	77	0.00
34	T	Trichloroethene	5.000	4.907	1.9	75	0.00
35	T	Methylcyclohexane	5.000	4.795	4.1	75	0.00
36	S	1,2-Dichloropropane-d6	5.000	5.767	-15.3	86	0.00
37	T	1,2-Dichloropropane	5.000	5.308	-6.2	80	0.00
38	T	Bromodichloromethane	5.000	5.185	-3.7	79	0.00
39	T	cis-1,3-Dichloropropene	5.000	5.081	-1.6	77	0.00
40	T	4-Methyl-2-pentanone	50.000	55.743	-11.5	84	0.00
41	S	Toluene-d8	5.000	5.546	-10.9	84	0.00
42	T	Toluene	5.000	4.898	2.0	76	0.00
43	S	trans-1,3-Dichloropropene-d	5.000	5.826	-16.5	88	0.00
44	T	trans-1,3-Dichloropropene	5.000	5.299	-6.0	80	0.00
45	T	1,1,2-Trichloroethane	5.000	5.151	-3.0	78	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA U\Data\VU010419\
Data File : VU029015.D
Acq On : 04 Jan 2019 09:37
Operator : MD/SY
Sample : VSTDCCC005
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_U
LabSampleId : VSTD00551

Quant Time: Jan 04 22:14:04 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR122818WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Thu Jan 03 03:47:56 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)
46	S	2-Hexanone-d5	50.000	61.872	-23.7	90	0.00
47	T	Tetrachloroethene	5.000	4.818	3.6	76	0.00
48	T	2-Hexanone	50.000	58.358	-16.7	86	0.00
49	T	Dibromochloromethane	5.000	5.283	-5.7	80	0.00
50	T	1,2-Dibromoethane	5.000	5.104	-2.1	76	0.00
51	T	Chlorobenzene	5.000	4.929	1.4	75	0.00
52	T	Ethylbenzene	5.000	4.860	2.8	75	0.00
53	T	m,p-xylene	5.000	4.924	1.5	74	0.00
54	T	o-xylene	5.000	4.887	2.3	74	0.00
55	T	Styrene	5.000	5.024	-0.5	76	0.00
56	T	Isopropylbenzene	5.000	4.818	3.6	74	0.00
57	S	1,1,2,2-Tetrachloroethane-d	5.000	5.537	-10.7	83	0.00
58	T	1,1,2,2-Tetrachloroethane	5.000	5.528	-10.6	82	0.00
59		1,2,3-Trichloropropane	5.000	5.352	-7.0	80	0.00
60	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	68	0.00
61	T	Bromoform	5.000	5.753	-15.1	81	0.00
62	T	1,3-Dichlorobenzene	5.000	4.899	2.0	72	0.00
63	T	1,4-Dichlorobenzene	5.000	4.859	2.8	73	0.00
64	S	1,2-Dichlorobenzene-d4	5.000	5.312	-6.2	77	0.00
65	T	1,2-Dichlorobenzene	5.000	5.066	-1.3	74	0.00
66	T	1,2-Dibromo-3-chloropropane	5.000	6.104	-22.1	88	0.00
67		1,3,5-Trichlorobenzene	5.000	4.959	0.8	71	0.00
68	T	1,2,4-trichlorobenzene	5.000	4.932	1.4	70	0.00
69		Naphthalene	5.000	5.446	-8.9	75	0.00
70	T	1,2,3-Trichlorobenzene	5.000	5.052	-1.0	74	0.00

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