Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV072018\
Data File : VV006657.D
Acq On : 20 Jul 2018 10:01
Operator : SY/MD
Sample : VSTDCCC005
Misc : 25 mL/MSVOA_V/WATER
ALS Vial : 2 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00571

Quant Time: Jul 26 02:03:47 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071918WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Sun Jul 22 23:06:02 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	1,4-Difluorobenzene	5.000	5.000	0.0	100	0.00
2	T	Dichlorodifluoromethane	5.000	5.592	-11.8	104	0.00
3	T	Chloromethane	5.000	5.400	-8.0	102	0.00
4	S	Vinyl Chloride-d3	5.000	4.945	1.1	94	0.00
5	T	Vinyl chloride	5.000	5.339	-6.8	101	0.00
6	T	Bromomethane	5.000	4.995	0.1	98	0.00
7	S	Chloroethane-d5	5.000	5.546	-10.9	100	0.00
8	T	Chloroethane	5.000	5.467	-9.3	105	0.00
9	T	Trichlorofluoromethane	5.000	5.621	-12.4	106	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	5.000	5.656	-13.1	108	0.00
11	S	1,1-Dichloroethene-d2	5.000	5.215	-4.3	101	0.00
12	T	1,1-Dichloroethene	5.000	5.513	-10.3	106	0.00
13	T	Acetone	50.000	50.832	-1.7	97	0.00
14	T	Carbon disulfide	5.000	5.502	-10.0	106	0.00
15	T	Methyl Acetate	5.000	5.012	-0.2	97	0.00
16	T	Methylene chloride	5.000	4.857	2.9	105	0.00
17	T	Methyl tert-butyl Ether	5.000	5.311	-6.2	101	0.00
18	T	trans-1,2-Dichloroethene	5.000	5.484	-9.7	104	0.00
19	T	1,1-Dichloroethane	5.000	5.684	-13.7	108	0.00
20	S	2-Butanone-d5	50.000	33.135	33.7#	63	0.00
21	T	2-Butanone	50.000	52.244	-4.5	98	0.00
22	T	cis-1,2-Dichloroethene	5.000	5.546	-10.9	105	0.00
23	T	Bromochloromethane	5.000	5.482	-9.6	104	0.00
24	S	Chloroform-d	5.000	5.178	-3.6	98	0.00
25	T	Chloroform	5.000	5.706	-14.1	108	0.00
26	S	1,2-Dichloroethane-d4	5.000	5.134	-2.7	97	0.00
27	T	1,2-Dichloroethane	5.000	5.515	-10.3	104	0.00
28	I	Chlorobenzene-d5	5.000	5.000	0.0	101	0.00
29	T	1,1,1-Trichloroethane	5.000	5.643	-12.9	107	0.00
30	T	Cyclohexane	5.000	5.609	-12.2	105	0.00
31	T	Carbon tetrachloride	5.000	5.624	-12.5	106	0.00
32	S	Benzene-d6	5.000	5.284	-5.7	99	0.00
33	T	Benzene	5.000	5.662	-13.2	106	0.00
34	T	Trichloroethene	5.000	5.474	-9.5	106	0.00
35	T	Methylcyclohexane	5.000	5.681	-13.6	106	0.00
36	S	1,2-Dichloropropane-d6	5.000	5.302	-6.0	101	0.00
37	T	1,2-Dichloropropane	5.000	5.643	-12.9	105	0.00
38	T	Bromodichloromethane	5.000	5.501	-10.0	104	0.00
39	T	cis-1,3-Dichloropropene	5.000	5.596	-11.9	104	0.00
40	T	4-Methyl-2-pentanone	50.000	52.203	-4.4	95	0.00
41	S	Toluene-d8	5.000	5.363	-7.3	100	0.00
42	T	Toluene	5.000	5.713	-14.3	106	0.00
43	S	trans-1,3-Dichloropropene-d	5.000	5.261	-5.2	99	0.00
44	T	trans-1,3-Dichloropropene	5.000	5.468	-9.4	101	0.00
45	T	1,1,2-Trichloroethane	5.000	5.382	-7.6	102	0.00

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		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
46	S	2-Hexanone-d5	50.000	47.007	6.0	86	0.00
47	T	Tetrachloroethene	5.000	5.638	-12.8	107	0.00
48	T	2-Hexanone	50.000	53.302	-6.6	96	0.00
49	T	Dibromochloromethane	5.000	5.574	-11.5	104	0.00
50	T	1,2-Dibromoethane	5.000	5.311	-6.2	99	0.00
51	T	Chlorobenzene	5.000	5.608	-12.2	106	0.00
52	T	Ethylbenzene	5.000	5.739	-14.8	106	0.00
53	T	m,p-xylene	5.000	5.855	-17.1	108	0.00
54	T	o-xylene	5.000	5.802	-16.0	107	0.00
55	T	Styrene	5.000	5.885	-17.7	106	0.00
56	T	Isopropylbenzene	5.000	5.820	-16.4	106	0.00
57	S	1,1,2,2-Tetrachloroethane-d	5.000	5.090	-1.8	96	0.00
58	T	1,1,2,2-Tetrachloroethane	5.000	5.412	-8.2	101	0.00
59		1,2,3-Trichloropropane	5.000	5.251	-5.0	100	0.00
60	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	101	0.00
61	T	Bromoform	5.000	5.237	-4.7	101	0.00
62	T	1,3-Dichlorobenzene	5.000	5.574	-11.5	106	0.00
63	T	1,4-Dichlorobenzene	5.000	5.535	-10.7	106	0.00
64	S	1,2-Dichlorobenzene-d4	5.000	5.282	-5.6	102	0.00
65	T	1,2-Dichlorobenzene	5.000	5.559	-11.2	105	0.00
66	T	1,2-Dibromo-3-chloropropane	5.000	5.191	-3.8	97	0.00
67		1,3,5-Trichlorobenzene	5.000	5.689	-13.8	108	0.00
68	T	1,2,4-trichlorobenzene	5.000	5.554	-11.1	104	0.00
69		Naphthalene	5.000	5.131	-2.6	95	0.00
70	T	1,2,3-Trichlorobenzene	5.000	5.505	-10.1	101	0.00

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