Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV100418\
Data File : VV008086.D
Acq On : 05 Oct 2018 02:31
Operator : SY/MD
Sample : VSTDCCC005EC
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00572

Quant Time: Oct 05 06:00:30 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Fri Oct 05 05:56:33 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	89	0.00
2	T	Dichlorodifluoromethane	5.000	5.017	-0.3	91	0.00
3	T	Chloromethane	5.000	4.534	9.3	81	0.00
4	S	Vinyl Chloride-d3	5.000	4.189	16.2	76	0.00
5	T	Vinyl chloride	5.000	5.037	-0.7	90	0.00
6	T	Bromomethane	5.000	2.047	59.1#	37	0.00
7	S	Chloroethane-d5	5.000	4.554	8.9	83	0.00
8	T	Chloroethane	5.000	5.234	-4.7	95	0.00
9	T	Trichlorofluoromethane	5.000	5.123	-2.5	92	0.00
10	T	1,1,2-Trichloro-1,2,2-trifl	5.000	5.277	-5.5	95	0.00
11	S	1,1-Dichloroethene-d2	5.000	4.418	11.6	81	0.00
12	T	1,1-Dichloroethene	5.000	5.109	-2.2	92	0.00
13	T	Acetone	50.000	47.836	4.3	92	0.00
14	T	Carbon disulfide	5.000	4.722	5.6	89	0.00
15	T	Methyl Acetate	5.000	5.003	-0.1	103	0.00
16	T	Methylene chloride	5.000	4.984	0.3	96	0.00
17	T	Methyl tert-butyl Ether	5.000	5.222	-4.4	97	0.00
18	T	trans-1,2-Dichloroethene	5.000	5.188	-3.8	95	0.00
19	T	1,1-Dichloroethane	5.000	5.576	-11.5	98	0.00
20	S	2-Butanone-d5	50.000	56.031	-12.1	107	0.00
21	T	2-Butanone	50.000	52.961	-5.9	102	0.00
22	T	cis-1,2-Dichloroethene	5.000	5.615	-12.3	101	0.00
23	T	Bromochloromethane	5.000	5.831	-16.6	105	0.00
24	S	Chloroform-d	5.000	5.312	-6.2	96	0.00
25	T	Chloroform	5.000	5.745	-14.9	103	0.00
26	S	1,2-Dichloroethane-d4	5.000	4.733	5.3	88	0.00
27	T	1,2-Dichloroethane	5.000	5.358	-7.2	96	0.00
28	I	Chlorobenzene-d5	5.000	5.000	0.0	87	0.00
29	T	1,1,1-Trichloroethane	5.000	5.780	-15.6	102	0.00
30	T	Cyclohexane	5.000	5.068	-1.4	90	0.00
31	T	Carbon tetrachloride	5.000	5.876	-17.5	103	0.00
32	S	Benzene-d6	5.000	5.105	-2.1	91	0.00
33	T	Benzene	5.000	5.740	-14.8	101	0.00
34	T	Trichloroethene	5.000	5.826	-16.5	105	0.00
35	T	Methylcyclohexane	5.000	4.981	0.4	88	0.00
36	S	1,2-Dichloropropane-d6	5.000	5.348	-7.0	96	0.00
37	T	1,2-Dichloropropane	5.000	5.988	-19.8	108	0.00
38	T	Bromodichloromethane	5.000	6.145	-22.9	109	0.00
39	T	cis-1,3-Dichloropropene	5.000	5.490	-9.8	98	0.00
40	T	4-Methyl-2-pentanone	50.000	55.945	-11.9	102	0.00
41	S	Toluene-d8	5.000	4.942	1.2	86	0.00
42	T	Toluene	5.000	5.771	-15.4	101	0.00
43	S	trans-1,3-Dichloropropene-d	5.000	4.783	4.3	89	0.00
44	T	trans-1,3-Dichloropropene	5.000	5.510	-10.2	98	0.00
45	T	1,1,2-Trichloroethane	5.000	5.796	-15.9	106	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV100418\
Data File : VV008086.D
Acq On : 05 Oct 2018 02:31
Operator : SY/MD
Sample : VSTDCCC005EC
Misc : 25.0 mL/MSVOA_V/WATER
ALS Vial : 33 Sample Multiplier: 1

Instrument : MSVOA_V
LabSampleId : VSTD00572

Quant Time: Oct 05 06:00:30 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR092718WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Fri Oct 05 05:56:33 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)
46	S	2-Hexanone-d5	50.000	54.245	-8.5	100	0.00
47	T	Tetrachloroethene	5.000	5.331	-6.6	95	0.00
48	T	2-Hexanone	50.000	57.820	-15.6	103	0.00
49	T	Dibromochloromethane	5.000	6.328	-26.6	113	0.00
50	T	1,2-Dibromoethane	5.000	5.845	-16.9	102	0.00
51	T	Chlorobenzene	5.000	5.575	-11.5	100	0.00
52	T	Ethylbenzene	5.000	5.461	-9.2	96	0.00
53	T	m,p-xylene	5.000	5.584	-11.7	99	0.00
54	T	o-xylene	5.000	5.549	-11.0	98	0.00
55	T	Styrene	5.000	5.716	-14.3	99	0.00
56	T	Isopropylbenzene	5.000	5.602	-12.0	97	0.00
57	S	1,1,2,2-Tetrachloroethane-d	5.000	5.489	-9.8	100	0.00
58	T	1,1,2,2-Tetrachloroethane	5.000	5.932	-18.6	107	0.00
59		1,2,3-Trichloropropane	5.000	5.713	-14.3	103	0.00
60	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	88	0.00
61	T	Bromoform	5.000	6.186	-23.7	118	0.00
62	T	1,3-Dichlorobenzene	5.000	5.389	-7.8	99	0.00
63	T	1,4-Dichlorobenzene	5.000	5.285	-5.7	99	0.00
64	S	1,2-Dichlorobenzene-d4	5.000	4.638	7.2	90	0.00
65	T	1,2-Dichlorobenzene	5.000	5.494	-9.9	101	0.00
66	T	1,2-Dibromo-3-chloropropane	5.000	5.662	-13.2	103	0.00
67		1,3,5-Trichlorobenzene	5.000	5.316	-6.3	98	0.00
68	T	1,2,4-trichlorobenzene	5.000	4.967	0.7	94	0.00
69		Naphthalene	5.000	4.576	8.5	85	0.00
70	T	1,2,3-Trichlorobenzene	5.000	5.249	-5.0	96	0.00

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