Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060716\
Data File : BM005705.D
Acq On : 07 Jun 2016 22:51
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02049

Quant Time: Jun 08 01:27:09 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 08 01:24:32 2016
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	148	0.00
2	S	1,4-Dioxane-d8	0.476	0.363	23.7#	116	0.00
3	S	Phenol-d5	1.758	1.744	0.8	146	0.01
4	S	Bis-(2-Chloroethyl)ether-d8	1.070	1.024	4.3	139	0.00
5	S	2-Chlorophenol-d4	1.451	1.456	-0.3	150#	0.00
6	S	4-Methylphenol-d8	1.391	1.413	-1.6	151#	0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	151#	0.00
8	S	Nitrobenzene-d5	0.154	0.156	-1.3	154#	0.01
9	S	2-Nitrophenol-d4	0.164	0.168	-2.4	154#	0.01
10	S	2,4-Dichlorophenol-d3	0.292	0.300	-2.7	156#	0.01
11		Naphthalene	1.017	1.001	1.6	150	0.01
12	S	4-Chloroaniline-d4	0.347	0.402	-15.9	157#	0.01
13		2-Methylnaphthalene	0.707	0.710	-0.4	153#	0.00
14		1-Methylnaphthalene	0.695	0.690	0.7	151#	0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
16	S	Dimethylphthalate-d6	1.538	1.580	-2.7	156#	0.00
17	S	Acenaphthylene-d8	2.000	2.036	-1.8	153#	0.01
18		Acenaphthylene	2.069	2.063	0.3	149	0.00
19	C	Acenaphthene	1.343	1.351	-0.6	152#	0.00
20	S	4-Nitrophenol-d4	0.220	0.218	0.9	153#	0.02
21	S	Fluorene-d10	1.321	1.310	0.8	150#	0.00
22		Fluorene	1.453	1.428	1.7	146	0.01
23	I	Phenanthrene-d10	1.000	1.000	0.0	142	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.094	0.033	64.9#	53	0.02
25		Phenanthrene	1.104	1.091	1.2	140	0.00
26	S	Anthracene-d10	0.950	0.949	0.1	141	0.01
27		Anthracene	1.122	1.113	0.8	140	0.01
28	C	Fluoranthene	1.128	0.984	12.8	122	0.01
29	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
30	S	Pyrene-d10	1.005	1.119	-11.3	112	0.01
31		Pyrene	1.262	1.407	-11.5	112	0.01
32		Benzo(a)anthracene	1.173	1.172	0.1	99	0.01
33		Chrysene	1.108	1.110	-0.2	101	0.00
34	I	Perylene-d12	1.000	1.000	0.0	108	0.01
35		Benzo(b)fluoranthene	1.206	1.132	6.1	103	0.01
36		Benzo(k)fluoranthene	1.132	1.112	1.8	106	0.01
37	S	Benzo(a)pyrene-d12	0.920	0.914	0.7	108	0.01
38	C	Benzo(a)pyrene	1.161	1.128	2.8	105	0.01
39		Indeno(1,2,3-cd)pyrene	1.374	1.355	1.4	106	0.02
40		Dibenzo(a,h)anthracene	1.153	1.130	2.0	105	0.03
41		Benzo(g,h,i)perylene	1.152	1.134	1.6	108	0.03

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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