Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM063016\
Data File : BM006163.D
Acq On : 30 Jun 2016 18:15
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02055

Quant Time: Jun 30 18:47:43 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM062816.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 29 00:57:09 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	139	0.00
2	S	1,4-Dioxane-d8	0.431	0.334	22.5#	135	0.00
3	S	Phenol-d5	1.634	1.385	15.2	142	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.963	0.840	12.8	141	0.00
5	S	2-Chlorophenol-d4	1.254	1.030	17.9	140	0.00
6	S	4-Methylphenol-d8	1.315	1.148	12.7	145	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	141	0.00
8	S	Nitrobenzene-d5	0.135	0.112	17.0	147	0.00
9	S	2-Nitrophenol-d4	0.154	0.128	16.9	145	0.00
10	S	2,4-Dichlorophenol-d3	0.281	0.232	17.4	144	0.00
11		Naphthalene	0.876	0.718	18.0	143	0.00
12	S	4-Chloroaniline-d4	0.222	0.178	19.8	124	0.00
13		2-Methylnaphthalene	0.665	0.553	16.8	144	0.00
14		1-Methylnaphthalene	0.632	0.526	16.8	146	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	141	0.00
16	S	Dimethylphthalate-d6	1.434	1.173	18.2	144	0.00
17	S	Acenaphthylene-d8	1.740	1.427	18.0	143	0.00
18		Acenaphthylene	1.822	1.497	17.8	141	0.00
19	C	Acenaphthene	1.162	0.957	17.6	141	0.00
20	S	4-Nitrophenol-d4	0.272	0.231	15.1	141	0.01
21	S	Fluorene-d10	1.223	1.000	18.2	143	0.00
22		Fluorene	1.323	1.092	17.5	141	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.071	0.087	-22.5#	256#	0.01
25		Phenanthrene	0.946	0.768	18.8	139	0.00
26	S	Anthracene-d10	0.802	0.666	17.0	141	0.00
27		Anthracene	0.973	0.803	17.5	140	0.00
28	C	Fluoranthene	1.159	0.947	18.3	136	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	123	0.00
30	S	Pyrene-d10	0.805	0.698	13.3	135	0.00
31		Pyrene	1.055	0.911	13.6	134	0.00
32		Benzo(a)anthracene	1.040	0.853	18.0	124	0.00
33		Chrysene	0.957	0.794	17.0	124	0.00
34	I	Perylene-d12	1.000	1.000	0.0	115	0.01
35		Benzo(b)fluoranthene	1.038	0.868	16.4	122	0.01
36		Benzo(k)fluoranthene	0.974	0.779	20.0#	113	0.00
37	S	Benzo(a)pyrene-d12	0.796	0.647	18.7	117	0.01
38	C	Benzo(a)pyrene	0.981	0.804	18.0	116	0.01
39		Indeno(1,2,3-cd)pyrene	1.019	0.849	16.7	112	0.01
40		Dibenzo(a,h)anthracene	0.843	0.710	15.8	113	0.02
41		Benzo(g,h,i)perylene	0.853	0.714	16.3	114	0.03

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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