Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\bm100615\
Data File : BM002854.D
Acq On : 06 Oct 2015 14:42
Operator : UM/IZ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Oct 06 15:31:19 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM100515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 06 15:21:07 2015
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	104	-0.02
2	SURR	1,4-Dioxane-d8	0.422	0.420	0.5	104	0.00
3		1,4-Dioxane	0.487	0.481	1.2	104	0.00
4	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
5		Naphthalene	1.063	1.061	0.2	105	0.00
6	SURR	2-Methylnaphthalene-d10	0.592	0.574	3.0	105	0.00
7		2-Methylnaphthalene	0.725	0.729	-0.6	106	0.00
8	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
9		Acenaphthylene	1.510	1.532	-1.5	111	0.00
10	C	Acenaphthene	1.516	1.515	0.1	107	0.00
11		Fluorene	1.709	1.707	0.1	107	-0.02
12	I	Phenanthrene-d10	1.000	1.000	0.0	107	-0.02
13		Pentachlorophenol	0.025	0.025	0.0	120	-0.02
14		Phenanthrene	1.227	1.234	-0.6	107	0.00
15		Anthracene	1.179	1.177	0.2	107	0.00
16	SURR	Fluoranthene-d10	0.989	0.997	-0.8	107	0.00
17	C	Fluoranthene	1.383	1.397	-1.0	107	0.00
18	I	Chrysene-d12	1.000	1.000	0.0	107	-0.01
19		Pyrene	1.586	1.574	0.8	107	0.00
20		Benzo(a)anthracene	1.336	1.347	-0.8	109	-0.01
21		Chrysene	1.309	1.319	-0.8	108	0.00
22	I	Perylene-d12	1.000	1.000	0.0	110	0.00
23		Benzo(b)fluoranthene	1.473	1.551	-5.3	114	0.00
24		Benzo(k)fluoranthene	1.431	1.332	6.9	105	0.00
25	C	Benzo(a)pyrene	1.340	1.348	-0.6	110	0.00
26		Indeno(1,2,3-cd)pyrene	1.364	1.355	0.7	111	0.00
27		Dibenzo(a,h)anthracene	1.080	1.075	0.5	112	-0.01
28		Benzo(g,h,i)perylene	1.176	1.166	0.9	111	-0.01

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