Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102114\
Data File : BE088178.D
Acq On : 22 Oct 2014 12:29
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Oct 22 13:03:11 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102014.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 21 02:11:14 2014
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	74	0.00
2		1,4-Dioxane	1.000	0.989	1.1	70	-0.08
3	I	Naphthalene-d8	5.000	5.000	0.0	79	0.00
4	S	Nitrobenzene-d5	1.000	1.355	-35.5#	103	0.00
5		Naphthalene	1.000	0.993	0.7	75	0.00
6		2-Methylnaphthalene	1.000	1.007	-0.7	76	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	89	0.00
8	S	2-Fluorobiphenyl	1.000	2.024	-102.4#	173	0.00
9		Acenaphthylene	1.000	0.906	9.4	77	0.01
10		Acenaphthene	1.000	0.928	7.2	82	0.00
11		Fluorene	1.000	0.961	3.9	82	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	87	0.00
13		Phenanthrene	1.000	1.442	-44.2#	102	0.00
14		Anthracene	1.000	0.991	0.9	85	0.00
15		Fluoranthene	1.000	1.534	-53.4#	134	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	64	0.00
17		Pyrene	1.000	1.796	-79.6#	116	0.00
18	S	Terphenyl-d14	1.000	2.523	-152.3#	160	0.00
19		Benzo(a)anthracene	1.000	1.253	-25.3#	83	0.00
20		Chrysene	1.000	1.300	-30.0#	84	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	1.166	-16.6	75	0.00
22	I	Perylene-d12	5.000	5.000	0.0	64	0.00
23		Benzo(b)fluoranthene	1.000	1.303	-30.3#	83	0.00
24		Benzo(k)fluoranthene	1.000	1.036	-3.6	67	0.00
25	C	Benzo(a)pyrene	1.000	1.228	-22.8#	77	0.00
26		Dibenzo(a,h)anthracene	1.000	0.894	10.6	58	0.00
27		Benzo(g,h,i)perylene	1.000	1.156	-15.6	74	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1