Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\Be102414\
Data File : BE088241.D
Acq On : 25 Oct 2014 10:04
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 27 15:06:17 2014
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-PAH-BE102414.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 27 13:04:40 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	77	0.00
2		1,4-Dioxane	1.000	1.024	-2.4	74	-0.09
3	I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
4	S	Nitrobenzene-d5	1.000	6.272	-527.2#	575	0.00
5		Naphthalene	1.000	1.023	-2.3	83	0.00
6		2-Methylnaphthalene	1.000	1.069	-6.9	84	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
8	S	2-Fluorobiphenyl	1.000	7.305	-630.5#	650	0.00
9		Acenaphthylene	1.000	1.026	-2.6	89	0.00
10		Acenaphthene	1.000	0.964	3.6	86	0.00
11		Fluorene	1.000	1.040	-4.0	90	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
13		Phenanthrene	1.000	1.175	-17.5	112	0.00
14		Anthracene	1.000	1.072	-7.2	100	0.00
15		Fluoranthene	1.000	1.533	-53.3#	144	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	85	0.00
17		Pyrene	1.000	1.590	-59.0#	126	0.00
18	S	Terphenyl-d14	1.000	10.233	-923.3#	820	0.00
19		Benzo(a)anthracene	1.000	1.314	-31.4#	110	0.00
20		Chrysene	1.000	1.157	-15.7	96	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	1.246	-24.6	123	0.00
22	I	Perylene-d12	5.000	5.000	0.0	84	0.00
23		Benzo(b)fluoranthene	1.000	1.415	-41.5#	127	0.00
24		Benzo(k)fluoranthene	1.000	1.019	-1.9	81	0.00
25	C	Benzo(a)pyrene	1.000	1.299	-29.9#	117	0.00
26		Dibenzo(a,h)anthracene	1.000	1.178	-17.8	107	0.00
27		Benzo(g,h,i)perylene	1.000	1.298	-29.8#	106	0.00

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