Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102114\
Data File : Be088141.d
Acq On : 21 Oct 2014 14:10
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 22 04:36:16 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	98	0.00
2		1,4-Dioxane	1.000	1.039	-3.9	97	-0.01
3	I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
4	S	Nitrobenzene-d5	1.000	1.001	-0.1	94	0.00
5		Naphthalene	1.000	1.039	-3.9	98	0.00
6		2-Methylnaphthalene	1.000	1.040	-4.0	97	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
8	S	2-Fluorobiphenyl	1.000	1.049	-4.9	98	0.00
9		Acenaphthylene	1.000	1.048	-4.8	97	0.00
10		Acenaphthene	1.000	1.037	-3.7	100	0.00
11		Fluorene	1.000	1.038	-3.8	97	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
13		Phenanthrene	1.000	0.929	7.1	79	0.00
14		Anthracene	1.000	0.999	0.1	95	0.00
15		Fluoranthene	1.000	0.973	2.7	95	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	96	0.00
17		Pyrene	1.000	0.996	0.4	96	0.00
18	S	Terphenyl-d14	1.000	0.999	0.1	94	0.00
19		Benzo(a)anthracene	1.000	0.939	6.1	92	0.00
20		Chrysene	1.000	1.026	-2.6	98	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	0.877	12.3	84	0.00
22	I	Perylene-d12	5.000	5.000	0.0	93	0.00
23		Benzo(b)fluoranthene	1.000	0.919	8.1	85	0.00
24		Benzo(k)fluoranthene	1.000	1.010	-1.0	93	0.00
25	C	Benzo(a)pyrene	1.000	0.920	8.0	83	0.00
26		Dibenzo(a,h)anthracene	1.000	0.894	10.6	83	0.00
27		Benzo(g,h,i)perylene	1.000	0.938	6.2	86	0.00

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