Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE102214\
Data File : BE088180.D
Acq On : 22 Oct 2014 13:40
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 27 04:18:20 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	77	0.00
2		1,4-Dioxane	1.000	1.011	-1.1	74	-0.06
3	I	Naphthalene-d8	5.000	5.000	0.0	78	0.00
4	S	Nitrobenzene-d5	1.000	1.027	-2.7	77	0.00
5		Naphthalene	1.000	1.036	-3.6	77	0.00
6		2-Methylnaphthalene	1.000	1.048	-4.8	78	0.00
7	I	Acenaphthene-d10	5.000	5.000	0.0	78	0.00
8	S	2-Fluorobiphenyl	1.000	1.055	-5.5	79	0.00
9		Acenaphthylene	1.000	1.062	-6.2	80	0.01
10		Acenaphthene	1.000	1.022	-2.2	79	0.00
11		Fluorene	1.000	1.019	-1.9	77	0.00
12	I	Phenanthrene-d10	5.000	5.000	0.0	73	0.00
13		Phenanthrene	1.000	0.926	7.4	59	0.00
14		Anthracene	1.000	0.996	0.4	71	0.00
15		Fluoranthene	1.000	0.862	13.8	63	0.00
16	I	Chrysene-d12	5.000	5.000	0.0	60	0.00
17		Pyrene	1.000	1.065	-6.5	64	0.00
18	S	Terphenyl-d14	1.000	1.080	-8.0	64	0.00
19		Benzo(a)anthracene	1.000	0.946	5.4	58	0.00
20		Chrysene	1.000	1.018	-1.8	61	0.00
21		Indeno(1,2,3-cd)pyrene	1.000	1.025	-2.5	62	0.00
22	I	Perylene-d12	5.000	5.000	0.0	59	0.00
23		Benzo(b)fluoranthene	1.000	0.914	8.6	54	0.00
24		Benzo(k)fluoranthene	1.000	1.003	-0.3	59	0.00
25	C	Benzo(a)pyrene	1.000	0.959	4.1	55	0.00
26		Dibenzo(a,h)anthracene	1.000	0.988	1.2	58	0.00
27		Benzo(g,h,i)perylene	1.000	1.042	-4.2	62	0.00

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