Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
Data File : BE089664.D
Acq On : 24 Feb 2015 6:36
Operator : TP/JJ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Quant Time: Feb 24 07:45:25 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE022315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 24 03:46:57 2015
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	109	0.00
2		1,4-Dioxane	1.000	1.187	-18.7	104	0.00
3		n-Nitrosodimethylamine	1.000	1.173	-17.3	109	0.00
4	I	Naphthalene-d8	5.000	5.000	0.0	109	0.00
5	S	Nitrobenzene-d5	1.000	1.110	-11.0	105	0.00
6		Naphthalene	1.000	1.159	-15.9	110	0.00
7		2-Methylnaphthalene	1.000	1.208	-20.8	112	0.00
8	I	Acenaphthene-d10	5.000	5.000	0.0	111	0.00
9	S	2-Fluorobiphenyl	1.000	1.197	-19.7	112	0.00
10		Acenaphthylene	1.000	1.208	-20.8	114	0.00
11		Acenaphthene	1.000	1.162	-16.2	112	0.00
12		Fluorene	1.000	1.186	-18.6	112	0.00
13	I	Phenanthrene-d10	5.000	5.000	0.0	94	0.00
14		Hexachlorobenzene	1.000	1.194	-19.4	99	0.00
15		Pentachlorophenol	1.000	0.960	4.0	90	0.00
16		Phenanthrene	1.000	1.165	-16.5	95	0.00
17		Anthracene	1.000	1.210	-21.0	97	0.00
18		Fluoranthene	1.000	1.204	-20.4	107	0.00
19	I	Chrysene-d12	5.000	5.000	0.0	104	0.00
20		Pyrene	1.000	1.219	-21.9	108	0.00
21	S	Terphenyl-d14	1.000	1.638	-63.8#	105	0.00
22		Benzo(a)anthracene	1.000	1.100	-10.0	99	0.00
23		Chrysene	1.000	1.169	-16.9	106	0.00
24		Indeno(1,2,3-cd)pyrene	1.000	0.932	6.8	93	0.00
25	I	Perylene-d12	5.000	5.000	0.0	102	0.00
26		Benzo(b)fluoranthene	1.000	0.870	13.0	93	0.00
27		Benzo(k)fluoranthene	1.000	1.162	-16.2	100	0.00
28	C	Benzo(a)pyrene	1.000	0.933	6.7	93	0.00
29		Dibenzo(a,h)anthracene	1.000	0.930	7.0	93	0.00
30		Benzo(g,h,i)perylene	1.000	1.028	-2.8	100	0.00

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