Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090333.D
Acq On : 4 Aug 2015 5:40
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 07:51:04 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 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	60	0.00
2		1,4-Dioxane	1.000	1.332	-33.2#	65	-0.01
3		n-Nitrosodimethylamine	1.000	1.090	-9.0	65	0.00
4	S	2-Fluorophenol	1.000	1.204	-20.4	75	0.00
5	S	Phenol-d6	1.000	1.283	-28.3#	83	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	61	0.00
7	S	Nitrobenzene-d5	1.000	1.015	-1.5	65	0.00
8		Nitrobenzene	1.000	0.978	2.2	64	0.00
9		Naphthalene	1.000	1.111	-11.1	69	0.00
10		Hexachlorobutadiene	1.000	1.085	-8.5	65	0.00
11		2-Methylnaphthalene	1.000	1.114	-11.4	69	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	55	0.00
13	S	2,4,6-Tribromophenol	1.000	1.391	-39.1#	98	0.00
14	S	2-Fluorobiphenyl	1.000	1.250	-25.0	68	0.00
15		Acenaphthylene	1.000	1.221	-22.1	67	0.00
16		Acenaphthene	1.000	1.172	-17.2	65	0.00
17		Fluorene	1.000	1.151	-15.1	67	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	49	0.02
19		4-Bromophenyl-phenylether	1.000	1.366	-36.6#	65	0.00
20		Hexachlorobenzene	1.000	1.330	-33.0#	64	0.00
21		Pentachlorophenol	1.000	1.358	-35.8#	87	0.00
22		Phenanthrene	1.000	1.199	-19.9	61	0.00
23		Anthracene	1.000	1.214	-21.4	63	0.00
24		Fluoranthene	1.000	1.199	-19.9	61	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	45	0.00
26		Pyrene	1.000	1.382	-38.2#	60	0.00
27	S	Terphenyl-d14	1.000	1.338	-33.8#	58	0.00
28		Benzo(a)anthracene	1.000	1.134	-13.4	58	0.00
29		Chrysene	1.000	1.674	-67.4#	64	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.215	-21.5	69	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.785	-78.5#	113	0.00
32	I	Perylene-d12	5.000	5.000	0.0	48	0.00
33		Benzo(b)fluoranthene	1.000	1.427	-42.7#	92	0.00
34		Benzo(k)fluoranthene	1.000	1.585	-58.5#	71	0.00
35	C	Benzo(a)pyrene	1.000	1.476	-47.6#	90	0.00
36		Dibenzo(a,h)anthracene	1.000	1.605	-60.5#	112	0.00
37		Benzo(g,h,i)perylene	1.000	1.680	-68.0#	93	-0.01

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