Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091115\
Data File : BE090783.D
Acq On : 11 Sep 2015 11:55
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 11 23:56:57 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	116	0.00
2		1,4-Dioxane	1.000	0.788	21.2	90	0.00
3		n-Nitrosodimethylamine	1.000	0.786	21.4	90	0.00
4	S	2-Fluorophenol	1.000	0.777	22.3	88	0.00
5	S	Phenol-d6	1.000	0.835	16.5	102	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	129	0.00
7	S	Nitrobenzene-d5	1.000	0.807	19.3	104	0.00
8		Nitrobenzene	1.000	0.834	16.6	110	0.00
9		Naphthalene	1.000	0.858	14.2	111	-0.01
10		Hexachlorobutadiene	1.000	0.848	15.2	105	0.00
11		2-Methylnaphthalene	1.000	0.892	10.8	118	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	141	0.00
13	S	2,4,6-Tribromophenol	1.000	0.762	23.8	106	0.00
14	S	2-Fluorobiphenyl	1.000	0.822	17.8	117	0.00
15		Acenaphthylene	1.000	0.832	16.8	122	0.00
16		Acenaphthene	1.000	0.856	14.4	119	0.00
17		Fluorene	1.000	0.816	18.4	116	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	133	0.00
19		4-Bromophenyl-phenylether	1.000	0.823	17.7	114	0.00
20		Hexachlorobenzene	1.000	0.840	16.0	111	0.00
21		Pentachlorophenol	1.000	0.863	13.7	83	0.00
22		Phenanthrene	1.000	0.871	12.9	116	-0.02
23		Anthracene	1.000	0.848	15.2	117	0.00
24		Fluoranthene	1.000	0.781	21.9	112	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	124	0.00
26		Pyrene	1.000	0.898	10.2	118	0.00
27	S	Terphenyl-d14	1.000	0.866	13.4	112	0.00
28		Benzo(a)anthracene	1.000	0.874	12.6	110	0.00
29		Chrysene	1.000	0.907	9.3	110	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.723	27.7#	96	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.867	13.3	116	0.00
32	I	Perylene-d12	5.000	5.000	0.0	128	0.00
33		Benzo(b)fluoranthene	1.000	0.835	16.5	110	0.00
34		Benzo(k)fluoranthene	1.000	0.861	13.9	112	0.00
35	C	Benzo(a)pyrene	1.000	0.843	15.7	115	0.00
36		Dibenzo(a,h)anthracene	1.000	0.861	13.9	115	-0.01
37		Benzo(g,h,i)perylene	1.000	0.807	19.3	109	-0.02

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