Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091415\
Data File : BE090796.D
Acq On : 14 Sep 2015 10:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 15 07:53:12 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	109	0.00
2		1,4-Dioxane	1.000	0.846	15.4	89	0.00
3		n-Nitrosodimethylamine	1.000	0.857	14.3	92	0.00
4	S	2-Fluorophenol	1.000	0.875	12.5	94	0.00
5	S	Phenol-d6	1.000	0.907	9.3	104	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	122	0.00
7	S	Nitrobenzene-d5	1.000	0.856	14.4	105	0.00
8		Nitrobenzene	1.000	0.854	14.6	108	0.00
9		Naphthalene	1.000	0.876	12.4	108	0.00
10		Hexachlorobutadiene	1.000	0.849	15.1	100	0.00
11		2-Methylnaphthalene	1.000	0.938	6.2	119	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	138	0.00
13	S	2,4,6-Tribromophenol	1.000	0.870	13.0	123	0.00
14	S	2-Fluorobiphenyl	1.000	0.835	16.5	117	0.00
15		Acenaphthylene	1.000	0.859	14.1	124	0.00
16		Acenaphthene	1.000	0.872	12.8	119	0.00
17		Fluorene	1.000	0.873	12.7	122	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	135	0.00
19		4-Bromophenyl-phenylether	1.000	0.858	14.2	121	-0.02
20		Hexachlorobenzene	1.000	0.829	17.1	112	0.00
21		Pentachlorophenol	1.000	0.964	3.6	106	0.00
22		Phenanthrene	1.000	0.907	9.3	123	-0.02
23		Anthracene	1.000	0.917	8.3	129	0.00
24		Fluoranthene	1.000	0.931	6.9	136	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	143	0.00
26		Pyrene	1.000	1.032	-3.2	155	0.00
27	S	Terphenyl-d14	1.000	0.960	4.0	143	0.00
28		Benzo(a)anthracene	1.000	0.954	4.6	138	0.00
29		Chrysene	1.000	0.949	5.1	131	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.993	0.7	159	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.986	1.4	152	-0.01
32	I	Perylene-d12	5.000	5.000	0.0	155	0.00
33		Benzo(b)fluoranthene	1.000	0.890	11.0	142	0.00
34		Benzo(k)fluoranthene	1.000	0.918	8.2	144	0.00
35	C	Benzo(a)pyrene	1.000	0.939	6.1	156	0.00
36		Dibenzo(a,h)anthracene	1.000	0.921	7.9	150	-0.01
37		Benzo(g,h,i)perylene	1.000	0.876	12.4	143	-0.01

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