Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071015\
Data File : BE090081.D
Acq On : 10 Jul 2015 12:37
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 10 23:08:10 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	120	-0.02
2		1,4-Dioxane	1.000	0.893	10.7	111	-0.01
3		n-Nitrosodimethylamine	1.000	0.815	18.5	102	-0.02
4	S	2-Fluorophenol	1.000	0.920	8.0	111	-0.01
5	S	Phenol-d6	1.000	0.948	5.2	116	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	120	-0.02
7	S	Nitrobenzene-d5	1.000	0.876	12.4	107	-0.02
8		Nitrobenzene	1.000	0.869	13.1	108	-0.02
9		Naphthalene	1.000	0.923	7.7	113	-0.01
10		Hexachlorobutadiene	1.000	0.956	4.4	117	-0.01
11		2-Methylnaphthalene	1.000	0.869	13.1	106	-0.02
12	I	Acenaphthene-d10	5.000	5.000	0.0	109	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.837	16.3	93	0.00
14	S	2-Fluorobiphenyl	1.000	0.919	8.1	103	-0.02
15		Acenaphthylene	1.000	0.918	8.2	103	-0.01
16		Acenaphthene	1.000	0.929	7.1	102	-0.01
17		Fluorene	1.000	0.895	10.5	100	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	96	-0.02
19		4-Bromophenyl-phenylether	1.000	0.909	9.1	89	-0.02
20		Hexachlorobenzene	1.000	0.901	9.9	90	-0.02
21		Pentachlorophenol	1.000	4.060	-306.0#	526	0.00
22		Phenanthrene	1.000	0.872	12.8	86	-0.02
23		Anthracene	1.000	0.892	10.8	87	0.00
24		Fluoranthene	1.000	0.857	14.3	84	-0.02
25	I	Chrysene-d12	5.000	5.000	0.0	84	-0.01
26		Pyrene	1.000	0.992	0.8	85	-0.02
27	S	Terphenyl-d14	1.000	0.938	6.2	79	-0.01
28		Benzo(a)anthracene	1.000	0.923	7.7	79	-0.01
29		Chrysene	1.000	0.902	9.8	77	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.992	0.8	96	-0.02
31		Indeno(1,2,3-cd)pyrene	1.000	0.942	5.8	81	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	79	-0.02
33		Benzo(b)fluoranthene	1.000	1.009	-0.9	81	-0.02
34		Benzo(k)fluoranthene	1.000	1.014	-1.4	81	-0.02
35	C	Benzo(a)pyrene	1.000	1.016	-1.6	81	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.969	3.1	78	-0.03
37		Benzo(g,h,i)perylene	1.000	0.979	2.1	79	-0.04

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