Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089942.D
Acq On : 26 Jun 2015 13:09
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 29 04:05:41 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 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	102	0.00
2		1,4-Dioxane	1.000	0.985	1.5	106	0.00
3		n-Nitrosodimethylamine	1.000	0.869	13.1	97	0.00
4	S	2-Fluorophenol	1.000	0.868	13.2	94	0.00
5	S	Phenol-d6	1.000	0.862	13.8	95	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
7	S	Nitrobenzene-d5	1.000	0.893	10.7	103	0.00
8		Nitrobenzene	1.000	0.879	12.1	101	0.00
9		Naphthalene	1.000	0.926	7.4	103	0.00
10		Hexachlorobutadiene	1.000	0.943	5.7	104	0.00
11		2-Methylnaphthalene	1.000	0.901	9.9	102	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
13	S	2,4,6-Tribromophenol	1.000	0.860	14.0	94	0.00
14	S	2-Fluorobiphenyl	1.000	0.903	9.7	101	0.00
15		Acenaphthylene	1.000	0.925	7.5	100	0.00
16		Acenaphthene	1.000	0.939	6.1	100	0.00
17		Fluorene	1.000	0.992	0.8	100	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
19		4-Bromophenyl-phenylether	1.000	0.866	13.4	98	-0.02
20		Hexachlorobenzene	1.000	0.917	8.3	100	0.00
21		Pentachlorophenol	1.000	0.979	2.1	108	-0.02
22		Phenanthrene	1.000	0.903	9.7	99	-0.02
23		Anthracene	1.000	0.892	10.8	97	0.00
24		Fluoranthene	1.000	0.905	9.5	98	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
26		Pyrene	1.000	0.930	7.0	97	0.00
27	S	Terphenyl-d14	1.000	0.886	11.4	97	0.00
28		Benzo(a)anthracene	1.000	0.922	7.8	100	0.00
29		Chrysene	1.000	0.891	10.9	98	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.846	15.4	100	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.971	2.9	102	0.00
32	I	Perylene-d12	5.000	5.000	0.0	100	0.00
33		Benzo(b)fluoranthene	1.000	0.966	3.4	102	0.00
34		Benzo(k)fluoranthene	1.000	0.957	4.3	102	0.00
35	C	Benzo(a)pyrene	1.000	0.930	7.0	100	0.00
36		Dibenzo(a,h)anthracene	1.000	0.951	4.9	102	0.00
37		Benzo(g,h,i)perylene	1.000	0.944	5.6	101	0.00

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