Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
Data File : BE090062.D
Acq On : 8 Jul 2015 16:51
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 09 01:19:06 2015
Quant Method : Z:\HPCHEM1\BNA_M\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	138	-0.01
2		1,4-Dioxane	1.000	0.845	15.5	121	-0.01
3		n-Nitrosodimethylamine	1.000	0.824	17.6	118	-0.02
4	S	2-Fluorophenol	1.000	0.857	14.3	118	-0.01
5	S	Phenol-d6	1.000	0.901	9.9	126	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	140	-0.01
7	S	Nitrobenzene-d5	1.000	0.865	13.5	123	-0.01
8		Nitrobenzene	1.000	0.858	14.2	124	-0.01
9		Naphthalene	1.000	0.918	8.2	131	-0.01
10		Hexachlorobutadiene	1.000	0.935	6.5	133	-0.01
11		2-Methylnaphthalene	1.000	0.883	11.7	126	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	132	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.667	33.3#	90	0.00
14	S	2-Fluorobiphenyl	1.000	0.902	9.8	122	-0.01
15		Acenaphthylene	1.000	0.911	8.9	124	-0.01
16		Acenaphthene	1.000	0.928	7.2	124	-0.01
17		Fluorene	1.000	0.892	10.8	121	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	122	-0.02
19		4-Bromophenyl-phenylether	1.000	0.880	12.0	110	-0.02
20		Hexachlorobenzene	1.000	0.888	11.2	113	-0.02
21		Pentachlorophenol	1.000	0.638	36.2#	67	0.00
22		Phenanthrene	1.000	0.874	12.6	110	-0.02
23		Anthracene	1.000	0.891	10.9	111	-0.02
24		Fluoranthene	1.000	0.873	12.7	109	-0.02
25	I	Chrysene-d12	5.000	5.000	0.0	113	-0.01
26		Pyrene	1.000	0.959	4.1	111	-0.02
27	S	Terphenyl-d14	1.000	0.914	8.6	104	-0.02
28		Benzo(a)anthracene	1.000	0.894	10.6	104	-0.01
29		Chrysene	1.000	0.902	9.8	104	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.702	29.8#	75	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.879	12.1	101	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	104	-0.02
33		Benzo(b)fluoranthene	1.000	0.995	0.5	105	-0.02
34		Benzo(k)fluoranthene	1.000	1.010	-1.0	106	-0.02
35	C	Benzo(a)pyrene	1.000	0.982	1.8	103	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.929	7.1	99	-0.03
37		Benzo(g,h,i)perylene	1.000	0.927	7.3	98	-0.04

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