Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070615\
Data File : BE090056.D
Acq On : 6 Jul 2015 15:01
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 07 01:13:18 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	61	-0.01
2		1,4-Dioxane	1.000	0.930	7.0	59	-0.02
3		n-Nitrosodimethylamine	1.000	0.787	21.3	50	-0.02
4	S	2-Fluorophenol	1.000	0.882	11.8	54	-0.01
5	S	Phenol-d6	1.000	0.883	11.7	55	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	58	-0.01
7	S	Nitrobenzene-d5	1.000	0.853	14.7	50	-0.01
8		Nitrobenzene	1.000	0.816	18.4	49	0.00
9		Naphthalene	1.000	0.908	9.2	54	-0.01
10		Hexachlorobutadiene	1.000	0.953	4.7	56	-0.01
11		2-Methylnaphthalene	1.000	0.833	16.7	49	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	49	0.00
13	S	2,4,6-Tribromophenol	1.000	0.637	36.3#	32	0.00
14	S	2-Fluorobiphenyl	1.000	0.993	0.7	50	0.00
15		Acenaphthylene	1.000	0.902	9.8	45	0.00
16		Acenaphthene	1.000	0.886	11.4	44	-0.01
17		Fluorene	1.000	0.869	13.1	43	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	42	-0.02
19		4-Bromophenyl-phenylether	1.000	0.905	9.5	39	-0.02
20		Hexachlorobenzene	1.000	0.921	7.9	40	-0.02
21		Pentachlorophenol	1.000	2.945	-194.5#	162	0.00
22		Phenanthrene	1.000	0.854	14.6	37	-0.02
23		Anthracene	1.000	0.852	14.8	36	0.00
24		Fluoranthene	1.000	0.817	18.3	35	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	37	-0.02
26		Pyrene	1.000	0.932	6.8	35	-0.01
27	S	Terphenyl-d14	1.000	0.926	7.4	34	-0.01
28		Benzo(a)anthracene	1.000	0.864	13.6	33	-0.01
29		Chrysene	1.000	0.877	12.3	33	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.620	38.0#	19	-0.02
31		Indeno(1,2,3-cd)pyrene	1.000	0.905	9.5	34	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	35	-0.02
33		Benzo(b)fluoranthene	1.000	1.002	-0.2	35	-0.02
34		Benzo(k)fluoranthene	1.000	0.974	2.6	34	-0.02
35	C	Benzo(a)pyrene	1.000	0.972	2.8	34	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.936	6.4	33	-0.04
37		Benzo(g,h,i)perylene	1.000	0.940	6.0	33	-0.04

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