Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070915\
Data File : BE090076.D
Acq On : 9 Jul 2015 16:56
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 10 02:02:40 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	114	-0.02
2		1,4-Dioxane	1.000	0.856	14.4	102	-0.01
3		n-Nitrosodimethylamine	1.000	0.801	19.9	95	-0.01
4	S	2-Fluorophenol	1.000	0.974	2.6	112	-0.01
5	S	Phenol-d6	1.000	1.077	-7.7	125	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	124	-0.02
7	S	Nitrobenzene-d5	1.000	0.881	11.9	111	-0.02
8		Nitrobenzene	1.000	0.865	13.5	111	-0.01
9		Naphthalene	1.000	0.920	8.0	116	-0.01
10		Hexachlorobutadiene	1.000	0.936	6.4	118	-0.01
11		2-Methylnaphthalene	1.000	0.918	8.2	116	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	120	-0.01
13	S	2,4,6-Tribromophenol	1.000	0.877	12.3	108	0.00
14	S	2-Fluorobiphenyl	1.000	0.919	8.1	113	-0.01
15		Acenaphthylene	1.000	0.915	8.5	113	0.00
16		Acenaphthene	1.000	0.931	6.9	113	-0.01
17		Fluorene	1.000	0.892	10.8	110	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	107	-0.02
19		4-Bromophenyl-phenylether	1.000	0.914	8.6	100	-0.02
20		Hexachlorobenzene	1.000	0.914	8.6	102	-0.02
21		Pentachlorophenol	1.000	1.492	-49.2#	191	0.00
22		Phenanthrene	1.000	0.881	11.9	97	-0.02
23		Anthracene	1.000	0.891	10.9	97	-0.02
24		Fluoranthene	1.000	0.856	14.4	94	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	93	-0.01
26		Pyrene	1.000	0.997	0.3	94	-0.01
27	S	Terphenyl-d14	1.000	0.949	5.1	88	-0.01
28		Benzo(a)anthracene	1.000	0.930	7.0	88	-0.01
29		Chrysene	1.000	0.902	9.8	85	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.085	-8.5	121	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	0.962	3.8	91	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	88	-0.02
33		Benzo(b)fluoranthene	1.000	1.045	-4.5	93	-0.02
34		Benzo(k)fluoranthene	1.000	1.002	-0.2	89	-0.02
35	C	Benzo(a)pyrene	1.000	1.036	-3.6	92	-0.02
36		Dibenzo(a,h)anthracene	1.000	0.985	1.5	89	-0.03
37		Benzo(g,h,i)perylene	1.000	0.992	0.8	89	-0.04

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