Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
Data File : BE090049.D
Acq On : 1 Jul 2015 20:12
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 02 05:16:31 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	102	0.00
2		1,4-Dioxane	1.000	0.859	14.1	91	0.00
3		n-Nitrosodimethylamine	1.000	0.825	17.5	88	-0.01
4	S	2-Fluorophenol	1.000	0.974	2.6	99	0.00
5	S	Phenol-d6	1.000	0.988	1.2	102	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	106	-0.02
7	S	Nitrobenzene-d5	1.000	0.871	12.9	94	0.00
8		Nitrobenzene	1.000	0.867	13.3	95	0.00
9		Naphthalene	1.000	0.920	8.0	99	0.00
10		Hexachlorobutadiene	1.000	0.935	6.5	100	0.00
11		2-Methylnaphthalene	1.000	0.896	10.4	96	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	1.000	0.878	12.2	89	0.00
14	S	2-Fluorobiphenyl	1.000	0.918	8.2	94	-0.01
15		Acenaphthylene	1.000	0.938	6.2	96	0.00
16		Acenaphthene	1.000	0.951	4.9	96	-0.01
17		Fluorene	1.000	0.910	9.0	93	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	93	-0.02
19		4-Bromophenyl-phenylether	1.000	0.875	12.5	83	-0.02
20		Hexachlorobenzene	1.000	0.940	6.0	91	-0.02
21		Pentachlorophenol	1.000	0.770	23.0	68	0.00
22		Phenanthrene	1.000	0.901	9.9	86	0.00
23		Anthracene	1.000	0.916	8.4	87	0.00
24		Fluoranthene	1.000	0.918	8.2	87	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	90	-0.01
26		Pyrene	1.000	0.952	4.8	87	-0.01
27	S	Terphenyl-d14	1.000	0.921	7.9	83	-0.01
28		Benzo(a)anthracene	1.000	0.957	4.3	88	-0.01
29		Chrysene	1.000	0.909	9.1	83	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	0.995	0.5	104	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	1.001	-0.1	92	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	87	-0.02
33		Benzo(b)fluoranthene	1.000	1.090	-9.0	96	-0.01
34		Benzo(k)fluoranthene	1.000	1.008	-0.8	88	-0.01
35	C	Benzo(a)pyrene	1.000	1.070	-7.0	93	-0.02
36		Dibenzo(a,h)anthracene	1.000	1.008	-0.8	89	-0.03
37		Benzo(g,h,i)perylene	1.000	1.005	-0.5	88	-0.03

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