Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
Data File : BE090013.D
Acq On : 30 Jun 2015 16:35
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 01 05:16:33 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	90	-0.01
2		1,4-Dioxane	1.000	0.885	11.5	83	-0.01
3		n-Nitrosodimethylamine	1.000	0.885	11.5	83	-0.02
4	S	2-Fluorophenol	1.000	1.046	-4.6	95	-0.01
5	S	Phenol-d6	1.000	1.034	-3.4	95	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	92	-0.01
7	S	Nitrobenzene-d5	1.000	0.935	6.5	87	0.00
8		Nitrobenzene	1.000	0.924	7.6	87	0.00
9		Naphthalene	1.000	0.924	7.6	86	-0.01
10		Hexachlorobutadiene	1.000	0.943	5.7	88	0.00
11		2-Methylnaphthalene	1.000	0.896	10.4	84	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	85	0.00
13	S	2,4,6-Tribromophenol	1.000	1.171	-17.1	101	0.00
14	S	2-Fluorobiphenyl	1.000	0.929	7.1	81	0.00
15		Acenaphthylene	1.000	0.943	5.7	83	0.00
16		Acenaphthene	1.000	0.938	6.2	81	0.00
17		Fluorene	1.000	0.912	8.8	80	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	77	-0.02
19		4-Bromophenyl-phenylether	1.000	0.909	9.1	71	0.00
20		Hexachlorobenzene	1.000	0.931	6.9	74	-0.02
21		Pentachlorophenol	1.000	1.065	-6.5	89	0.00
22		Phenanthrene	1.000	0.911	8.9	72	0.00
23		Anthracene	1.000	0.980	2.0	77	0.00
24		Fluoranthene	1.000	0.965	3.5	75	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	80	0.00
26		Pyrene	1.000	0.944	5.6	77	0.00
27	S	Terphenyl-d14	1.000	0.915	8.5	73	0.00
28		Benzo(a)anthracene	1.000	0.965	3.5	79	0.00
29		Chrysene	1.000	0.905	9.5	73	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.393	-39.3#	146	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.137	-13.7	93	-0.02
32	I	Perylene-d12	5.000	5.000	0.0	83	0.00
33		Benzo(b)fluoranthene	1.000	1.066	-6.6	89	0.00
34		Benzo(k)fluoranthene	1.000	1.013	-1.3	84	0.00
35	C	Benzo(a)pyrene	1.000	1.090	-9.0	90	-0.01
36		Dibenzo(a,h)anthracene	1.000	1.086	-8.6	91	-0.01
37		Benzo(g,h,i)perylene	1.000	1.096	-9.6	92	-0.02

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