Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089962.D
Acq On : 27 Jun 2015 4:34
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jun 27 06:30:22 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 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	86	0.00
2		1,4-Dioxane	1.000	1.054	-5.4	95	0.00
3		n-Nitrosodimethylamine	1.000	0.828	17.2	78	0.00
4	S	2-Fluorophenol	1.000	1.362	-36.2#	124	0.00
5	S	Phenol-d6	1.000	2.304	-130.4#	213	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	93	0.00
7	S	Nitrobenzene-d5	1.000	1.395	-39.5#	145	0.00
8		Nitrobenzene	1.000	0.758	24.2	79	0.00
9		Naphthalene	1.000	0.830	17.0	83	0.00
10		Hexachlorobutadiene	1.000	0.856	14.4	85	0.00
11		2-Methylnaphthalene	1.000	0.788	21.2	80	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
13	S	2,4,6-Tribromophenol	1.000	1.721	-72.1#	221	0.00
14	S	2-Fluorobiphenyl	1.000	1.795	-79.5#	184	0.00
15		Acenaphthylene	1.000	0.785	21.5	77	0.00
16		Acenaphthene	1.000	0.808	19.2	78	0.00
17		Fluorene	1.000	0.829	17.1	76	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	86	0.00
19		4-Bromophenyl-phenylether	1.000	0.754	24.6	75	-0.02
20		Hexachlorobenzene	1.000	0.821	17.9	79	0.00
21		Pentachlorophenol	1.000	0.849	15.1	72	-0.02
22		Phenanthrene	1.000	0.781	21.9	75	-0.02
23		Anthracene	1.000	0.758	24.2	72	0.00
24		Fluoranthene	1.000	0.740	26.0#	70	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	81	0.00
26		Pyrene	1.000	0.823	17.7	70	0.00
27	S	Terphenyl-d14	1.000	1.935	-93.5#	173	0.00
28		Benzo(a)anthracene	1.000	0.763	23.7	68	0.00
29		Chrysene	1.000	0.758	24.2	68	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.757	24.3	73	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.745	25.5#	64	0.00
32	I	Perylene-d12	5.000	5.000	0.0	77	0.00
33		Benzo(b)fluoranthene	1.000	0.803	19.7	65	0.00
34		Benzo(k)fluoranthene	1.000	0.807	19.3	66	0.00
35	C	Benzo(a)pyrene	1.000	0.764	23.6#	63	0.00
36		Dibenzo(a,h)anthracene	1.000	0.780	22.0	65	0.00
37		Benzo(g,h,i)perylene	1.000	0.775	22.5	64	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1