Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080415\
Data File : BE090344.D
Acq On : 4 Aug 2015 15:34
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 04 17:41:38 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 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	95	-0.01
2		1,4-Dioxane	1.000	1.208	-20.8	94	-0.02
3		n-Nitrosodimethylamine	1.000	1.165	-16.5	109	-0.02
4	S	2-Fluorophenol	1.000	1.137	-13.7	111	-0.01
5	S	Phenol-d6	1.000	1.050	-5.0	108	-0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	108	-0.02
7	S	Nitrobenzene-d5	1.000	0.961	3.9	108	0.00
8		Nitrobenzene	1.000	0.964	3.6	110	0.00
9		Naphthalene	1.000	0.980	2.0	106	-0.02
10		Hexachlorobutadiene	1.000	0.915	8.5	96	0.00
11		2-Methylnaphthalene	1.000	0.976	2.4	106	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	107	-0.01
13	S	2,4,6-Tribromophenol	1.000	1.231	-23.1	164	-0.02
14	S	2-Fluorobiphenyl	1.000	1.028	-2.8	109	-0.01
15		Acenaphthylene	1.000	0.980	2.0	105	-0.01
16		Acenaphthene	1.000	0.961	3.9	104	-0.01
17		Fluorene	1.000	0.984	1.6	112	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
19		4-Bromophenyl-phenylether	1.000	0.978	2.2	106	-0.02
20		Hexachlorobenzene	1.000	0.954	4.6	104	0.00
21		Pentachlorophenol	1.000	10.710	-971.0#	3601	-0.03
22		Phenanthrene	1.000	0.978	2.2	114	-0.02
23		Anthracene	1.000	0.970	3.0	115	-0.02
24		Fluoranthene	1.000	1.003	-0.3	117	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	142	-0.01
26		Pyrene	1.000	0.884	11.6	120	-0.01
27	S	Terphenyl-d14	1.000	0.978	2.2	133	-0.02
28		Benzo(a)anthracene	1.000	1.017	-1.7	162	-0.01
29		Chrysene	1.000	0.959	4.1	118	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.275	-27.5#	230	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.639	-63.9#	322	-0.04
32	I	Perylene-d12	5.000	5.000	0.0	165	-0.02
33		Benzo(b)fluoranthene	1.000	1.143	-14.3	234	-0.02
34		Benzo(k)fluoranthene	1.000	0.863	13.7	133	-0.01
35	C	Benzo(a)pyrene	1.000	1.150	-15.0	223	-0.02
36		Dibenzo(a,h)anthracene	1.000	1.603	-60.3#	385	-0.05
37		Benzo(g,h,i)perylene	1.000	1.189	-18.9	212	-0.04

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