Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
Data File : BE090530.D
Acq On : 14 Aug 2015 21:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 17 04:19:26 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	105	0.00
2		1,4-Dioxane	1.000	1.132	-13.2	102	0.00
3		n-Nitrosodimethylamine	1.000	1.008	-0.8	107	0.00
4	S	2-Fluorophenol	1.000	1.166	-16.6	121	0.00
5	S	Phenol-d6	1.000	1.139	-13.9	124	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
7	S	Nitrobenzene-d5	1.000	1.102	-10.2	119	0.00
8		Nitrobenzene	1.000	1.106	-10.6	121	0.00
9		Naphthalene	1.000	1.004	-0.4	104	0.00
10		Hexachlorobutadiene	1.000	0.984	1.6	101	0.00
11		2-Methylnaphthalene	1.000	1.027	-2.7	108	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	104	0.00
13	S	2,4,6-Tribromophenol	1.000	1.225	-22.5	141	0.00
14	S	2-Fluorobiphenyl	1.000	1.005	-0.5	105	0.00
15		Acenaphthylene	1.000	1.026	-2.6	109	0.00
16		Acenaphthene	1.000	1.011	-1.1	107	0.00
17		Fluorene	1.000	0.998	0.2	105	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
19		4-Bromophenyl-phenylether	1.000	1.097	-9.7	111	0.00
20		Hexachlorobenzene	1.000	1.026	-2.6	102	0.00
21		Pentachlorophenol	1.000	1.303	-30.3#	192	0.00
22		Phenanthrene	1.000	0.995	0.5	100	0.00
23		Anthracene	1.000	1.071	-7.1	110	0.00
24		Fluoranthene	1.000	1.069	-6.9	109	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
26		Pyrene	1.000	1.059	-5.9	109	0.00
27	S	Terphenyl-d14	1.000	1.081	-8.1	110	0.00
28		Benzo(a)anthracene	1.000	1.028	-2.8	109	0.00
29		Chrysene	1.000	1.015	-1.5	103	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.459	-45.9#	180	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.115	-11.5	121	0.00
32	I	Perylene-d12	5.000	5.000	0.0	108	0.00
33		Benzo(b)fluoranthene	1.000	1.085	-8.5	120	0.00
34		Benzo(k)fluoranthene	1.000	1.014	-1.4	108	0.00
35	C	Benzo(a)pyrene	1.000	1.073	-7.3	120	0.00
36		Dibenzo(a,h)anthracene	1.000	1.096	-9.6	125	0.00
37		Benzo(g,h,i)perylene	1.000	1.083	-8.3	119	0.00

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