Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090427.D
Acq On : 11 Aug 2015 7:04
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 07:37:47 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 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	126	0.00
2		1,4-Dioxane	1.000	1.076	-7.6	120	0.00
3		n-Nitrosodimethylamine	1.000	1.068	-6.8	136	0.00
4	S	2-Fluorophenol	1.000	1.517	-51.7#	206	0.00
5	S	Phenol-d6	1.000	1.242	-24.2	183	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	133	0.00
7	S	Nitrobenzene-d5	1.000	1.042	-4.2	151	0.00
8		Nitrobenzene	1.000	1.017	-1.7	157	0.00
9		Naphthalene	1.000	1.027	-2.7	138	0.00
10		Hexachlorobutadiene	1.000	1.074	-7.4	131	0.00
11		2-Methylnaphthalene	1.000	1.130	-13.0	157	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	144	0.00
13	S	2,4,6-Tribromophenol	1.000	1.635	-63.5#	290	0.00
14	S	2-Fluorobiphenyl	1.000	1.055	-5.5	150	0.00
15		Acenaphthylene	1.000	1.082	-8.2	158	0.00
16		Acenaphthene	1.000	1.013	-1.3	146	0.00
17		Fluorene	1.000	1.016	-1.6	145	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	136	0.00
19		4-Bromophenyl-phenylether	1.000	1.025	-2.5	143	0.00
20		Hexachlorobenzene	1.000	0.925	7.5	125	0.00
21		Pentachlorophenol	1.000	1.193	-19.3	202	0.00
22		Phenanthrene	1.000	1.020	-2.0	140	0.00
23		Anthracene	1.000	1.121	-12.1	156	-0.02
24		Fluoranthene	1.000	1.116	-11.6	157	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	132	0.00
26		Pyrene	1.000	1.158	-15.8	159	0.00
27	S	Terphenyl-d14	1.000	1.173	-17.3	156	0.00
28		Benzo(a)anthracene	1.000	1.096	-9.6	161	0.00
29		Chrysene	1.000	1.034	-3.4	133	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.402	-40.2#	227	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.093	-9.3	181	0.00
32	I	Perylene-d12	5.000	5.000	0.0	141	0.00
33		Benzo(b)fluoranthene	1.000	1.112	-11.2	176	0.00
34		Benzo(k)fluoranthene	1.000	0.945	5.5	129	0.00
35	C	Benzo(a)pyrene	1.000	1.076	-7.6	167	0.00
36		Dibenzo(a,h)anthracene	1.000	1.092	-9.2	190	0.00
37		Benzo(g,h,i)perylene	1.000	1.112	-11.2	162	0.00

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