Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090416.D
Acq On : 11 Aug 2015 00:27
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 04:00:42 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	130	0.00
2		1,4-Dioxane	1.000	1.040	-4.0	120	0.00
3		n-Nitrosodimethylamine	1.000	1.066	-6.6	139	0.00
4	S	2-Fluorophenol	1.000	1.567	-56.7#	218	0.00
5	S	Phenol-d6	1.000	1.271	-27.1#	194	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	137	0.00
7	S	Nitrobenzene-d5	1.000	1.080	-8.0	162	0.00
8		Nitrobenzene	1.000	1.047	-4.7	166	0.00
9		Naphthalene	1.000	1.019	-1.9	140	0.00
10		Hexachlorobutadiene	1.000	1.063	-6.3	133	0.00
11		2-Methylnaphthalene	1.000	1.155	-15.5	165	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
13	S	2,4,6-Tribromophenol	1.000	1.786	-78.6#	331	0.00
14	S	2-Fluorobiphenyl	1.000	1.065	-6.5	155	0.00
15		Acenaphthylene	1.000	1.109	-10.9	165	0.00
16		Acenaphthene	1.000	1.011	-1.1	149	0.00
17		Fluorene	1.000	1.020	-2.0	149	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	139	0.00
19		4-Bromophenyl-phenylether	1.000	1.079	-7.9	153	0.00
20		Hexachlorobenzene	1.000	0.947	5.3	131	0.00
21		Pentachlorophenol	1.000	1.208	-20.8	210	0.00
22		Phenanthrene	1.000	1.047	-4.7	147	0.00
23		Anthracene	1.000	1.177	-17.7	168	-0.02
24		Fluoranthene	1.000	1.168	-16.8	168	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	141	0.00
26		Pyrene	1.000	1.167	-16.7	170	0.00
27	S	Terphenyl-d14	1.000	1.177	-17.7	167	0.00
28		Benzo(a)anthracene	1.000	1.099	-9.9	172	0.00
29		Chrysene	1.000	1.035	-3.5	142	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.577	-57.7#	280	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.200	-20.0	218	0.00
32	I	Perylene-d12	5.000	5.000	0.0	151	0.00
33		Benzo(b)fluoranthene	1.000	1.122	-12.2	191	0.00
34		Benzo(k)fluoranthene	1.000	1.012	-1.2	148	0.00
35	C	Benzo(a)pyrene	1.000	1.142	-14.2	194	0.00
36		Dibenzo(a,h)anthracene	1.000	1.197	-19.7	231	0.00
37		Benzo(g,h,i)perylene	1.000	1.207	-20.7	189	-0.01

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