Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081315\
Data File : BE090512.D
Acq On : 13 Aug 2015 22:28
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 14 04:35:45 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	138	0.00
2		1,4-Dioxane	1.000	0.815	18.5	103	0.00
3		n-Nitrosodimethylamine	1.000	0.939	6.1	130	0.00
4	S	2-Fluorophenol	1.000	1.779	-77.9#	263	0.00
5	S	Phenol-d6	1.000	1.401	-40.1#	232	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	155	0.00
7	S	Nitrobenzene-d5	1.000	0.980	2.0	163	0.00
8		Nitrobenzene	1.000	0.964	3.6	170	0.00
9		Naphthalene	1.000	0.904	9.6	141	0.00
10		Hexachlorobutadiene	1.000	0.799	20.1	116	0.00
11		2-Methylnaphthalene	1.000	0.986	1.4	160	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	155	0.00
13	S	2,4,6-Tribromophenol	1.000	1.819	-81.9#	355	0.00
14	S	2-Fluorobiphenyl	1.000	0.919	8.1	140	0.00
15		Acenaphthylene	1.000	0.989	1.1	155	0.00
16		Acenaphthene	1.000	0.967	3.3	150	0.00
17		Fluorene	1.000	0.908	9.2	139	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	144	0.00
19		4-Bromophenyl-phenylether	1.000	0.871	12.9	129	0.00
20		Hexachlorobenzene	1.000	0.765	23.5	110	0.00
21		Pentachlorophenol	1.000	1.646	-64.6#	328	0.00
22		Phenanthrene	1.000	0.888	11.2	129	0.00
23		Anthracene	1.000	1.013	-1.3	150	-0.02
24		Fluoranthene	1.000	0.972	2.8	146	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	123	0.00
26		Pyrene	1.000	1.150	-15.0	147	0.00
27	S	Terphenyl-d14	1.000	1.049	-4.9	130	-0.01
28		Benzo(a)anthracene	1.000	1.014	-1.4	139	0.00
29		Chrysene	1.000	0.839	16.1	101	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	2.079	-107.9#	346	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.097	-9.7	170	0.00
32	I	Perylene-d12	5.000	5.000	0.0	132	0.00
33		Benzo(b)fluoranthene	1.000	1.071	-7.1	158	0.00
34		Benzo(k)fluoranthene	1.000	0.837	16.3	107	0.00
35	C	Benzo(a)pyrene	1.000	1.072	-7.2	157	0.00
36		Dibenzo(a,h)anthracene	1.000	1.089	-8.9	178	-0.01
37		Benzo(g,h,i)perylene	1.000	1.099	-9.9	150	0.00

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