Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090445.D
Acq On : 12 Aug 2015 1:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 06:42:22 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	117	-0.01
2		1,4-Dioxane	1.000	1.080	-8.0	111	0.00
3		n-Nitrosodimethylamine	1.000	1.046	-4.6	123	0.00
4	S	2-Fluorophenol	1.000	1.362	-36.2#	171	0.00
5	S	Phenol-d6	1.000	1.192	-19.2	161	-0.01
6	I	Naphthalene-d8	5.000	5.000	0.0	123	0.00
7	S	Nitrobenzene-d5	1.000	1.013	-1.3	135	-0.01
8		Nitrobenzene	1.000	0.999	0.1	142	0.00
9		Naphthalene	1.000	1.011	-1.1	126	0.00
10		Hexachlorobutadiene	1.000	1.090	-9.0	123	0.00
11		2-Methylnaphthalene	1.000	1.086	-8.6	140	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	133	0.00
13	S	2,4,6-Tribromophenol	1.000	1.396	-39.6#	220	0.00
14	S	2-Fluorobiphenyl	1.000	1.033	-3.3	135	0.00
15		Acenaphthylene	1.000	1.042	-4.2	140	0.00
16		Acenaphthene	1.000	1.010	-1.0	134	0.00
17		Fluorene	1.000	1.021	-2.1	134	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	131	0.00
19		4-Bromophenyl-phenylether	1.000	0.986	1.4	132	0.00
20		Hexachlorobenzene	1.000	0.941	5.9	123	0.00
21		Pentachlorophenol	1.000	1.210	-21.0	199	0.00
22		Phenanthrene	1.000	0.990	1.0	131	0.00
23		Anthracene	1.000	1.077	-7.7	145	-0.02
24		Fluoranthene	1.000	1.034	-3.4	141	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	127	0.00
26		Pyrene	1.000	1.098	-9.8	144	0.00
27	S	Terphenyl-d14	1.000	1.099	-9.9	140	0.00
28		Benzo(a)anthracene	1.000	1.046	-4.6	147	0.00
29		Chrysene	1.000	1.014	-1.4	125	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.109	-10.9	164	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.042	-4.2	164	0.00
32	I	Perylene-d12	5.000	5.000	0.0	130	0.00
33		Benzo(b)fluoranthene	1.000	1.065	-6.5	154	0.00
34		Benzo(k)fluoranthene	1.000	0.953	4.7	120	0.00
35	C	Benzo(a)pyrene	1.000	1.036	-3.6	147	0.00
36		Dibenzo(a,h)anthracene	1.000	1.061	-6.1	169	0.00
37		Benzo(g,h,i)perylene	1.000	1.061	-6.1	143	0.00

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