Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060815\
Data File : BM001623.D
Acq On : 09 Jun 2015 20:11
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 10 04:04:22 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 10 03:26:36 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	107	0.00
2		1,4-Dioxane	1.000	0.934	6.6	105	0.00
3		n-Nitrosodimethylamine	1.000	0.980	2.0	106	0.00
4	S	2-Fluorophenol	1.000	1.002	-0.2	109	-0.02
5	S	Phenol-d6	1.000	1.013	-1.3	112	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	111	0.00
7	S	Nitrobenzene-d5	1.000	0.992	0.8	113	0.00
8		Nitrobenzene	1.000	0.991	0.9	114	0.00
9		Naphthalene	1.000	0.997	0.3	113	0.00
10		Hexachlorobutadiene	1.000	0.968	3.2	109	0.00
11		2-Methylnaphthalene	1.000	1.006	-0.6	115	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
13	S	2,4,6-Tribromophenol	1.000	0.957	4.3	122	0.00
14	S	2-Fluorobiphenyl	1.000	0.967	3.3	116	0.00
15		Acenaphthylene	1.000	0.982	1.8	120	0.00
16		Acenaphthene	1.000	0.975	2.5	118	0.00
17		Fluorene	1.000	1.127	-12.7	137	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	110	0.03
19		4-Bromophenyl-phenylether	1.000	1.055	-5.5	116	0.00
20		Hexachlorobenzene	1.000	1.293	-29.3#	147	0.00
21		Pentachlorophenol	1.000	1.239	-23.9	138	0.00
22		Phenanthrene	1.000	0.950	5.0	95	0.00
23		Anthracene	1.000	1.361	-36.1#	156	0.00
24		Fluoranthene	1.000	1.094	-9.4	126	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	128	-0.01
26		Pyrene	1.000	0.956	4.4	128	0.00
27	S	Terphenyl-d14	1.000	0.975	2.5	129	0.00
28		Benzo(a)anthracene	1.000	0.960	4.0	129	0.00
29		Chrysene	1.000	0.975	2.5	126	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.888	11.2	127	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.987	1.3	132	0.00
32	I	Perylene-d12	5.000	5.000	0.0	130	0.00
33		Benzo(b)fluoranthene	1.000	1.000	0.0	135	0.00
34		Benzo(k)fluoranthene	1.000	0.960	4.0	125	0.00
35	C	Benzo(a)pyrene	1.000	0.980	2.0	132	0.00
36		Dibenzo(a,h)anthracene	1.000	0.986	1.4	133	0.00
37		Benzo(g,h,i)perylene	1.000	0.971	2.9	131	0.00

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