Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
Data File : BM001538.D
Acq On : 03 Jun 2015 07:20
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 03 08:05:37 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 03 05:42:14 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	99	0.00
2		1,4-Dioxane	1.000	0.994	0.6	103	0.00
3		n-Nitrosodimethylamine	1.000	1.006	-0.6	97	0.00
4	S	2-Fluorophenol	1.000	1.017	-1.7	103	-0.02
5	S	Phenol-d6	1.000	1.010	-1.0	103	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
7	S	Nitrobenzene-d5	1.000	1.070	-7.0	108	0.00
8		Nitrobenzene	1.000	1.063	-6.3	107	-0.01
9		Naphthalene	1.000	0.979	2.1	98	0.00
10		Hexachlorobutadiene	1.000	0.985	1.5	98	0.00
11		2-Methylnaphthalene	1.000	0.983	1.7	97	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
13	S	2,4,6-Tribromophenol	1.000	1.366	-36.6#	138	0.00
14	S	2-Fluorobiphenyl	1.000	0.977	2.3	97	0.00
15		Acenaphthylene	1.000	0.991	0.9	99	0.00
16		Acenaphthene	1.000	0.995	0.5	97	0.00
17		Fluorene	1.000	0.777	22.3	76	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	100	0.00
19		4-Bromophenyl-phenylether	1.000	1.189	-18.9	121	0.00
20		Hexachlorobenzene	1.000	0.670	33.0#	67	0.00
21		Pentachlorophenol	1.000	0.857	14.3	85	0.00
22		Phenanthrene	1.000	1.259	-25.9#	137	0.00
23		Anthracene	1.000	0.705	29.5#	70	0.00
24		Fluoranthene	1.000	0.961	3.9	99	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	99	-0.01
26		Pyrene	1.000	0.990	1.0	100	0.00
27	S	Terphenyl-d14	1.000	0.980	2.0	96	0.00
28		Benzo(a)anthracene	1.000	1.041	-4.1	108	-0.01
29		Chrysene	1.000	0.996	0.4	101	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.016	-1.6	115	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	1.046	-4.6	106	-0.02
32	I	Perylene-d12	5.000	5.000	0.0	106	0.00
33		Benzo(b)fluoranthene	1.000	0.966	3.4	99	0.00
34		Benzo(k)fluoranthene	1.000	1.016	-1.6	112	0.00
35	C	Benzo(a)pyrene	1.000	1.010	-1.0	107	0.00
36		Dibenzo(a,h)anthracene	1.000	0.991	0.9	105	-0.02
37		Benzo(g,h,i)perylene	1.000	0.998	0.2	106	0.00

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