Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060815\
Data File : BM001612.D
Acq On : 09 Jun 2015 10:33
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 10 03:26:24 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 09 17:36:38 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	109	0.00
2		1,4-Dioxane	1.000	0.906	9.4	104	0.00
3		n-Nitrosodimethylamine	1.000	1.017	-1.7	112	0.00
4	S	2-Fluorophenol	1.000	0.972	2.8	108	-0.02
5	S	Phenol-d6	1.000	0.987	1.3	111	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	111	0.00
7	S	Nitrobenzene-d5	1.000	0.983	1.7	111	0.00
8		Nitrobenzene	1.000	0.983	1.7	112	0.00
9		Naphthalene	1.000	1.012	-1.2	114	0.00
10		Hexachlorobutadiene	1.000	0.991	0.9	111	0.00
11		2-Methylnaphthalene	1.000	1.011	-1.1	115	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	117	0.00
13	S	2,4,6-Tribromophenol	1.000	0.934	6.6	117	0.00
14	S	2-Fluorobiphenyl	1.000	0.990	1.0	116	0.00
15		Acenaphthylene	1.000	0.976	2.4	117	0.00
16		Acenaphthene	1.000	0.977	2.3	116	0.00
17		Fluorene	1.000	1.022	-2.2	122	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	110	0.03
19		4-Bromophenyl-phenylether	1.000	1.052	-5.2	116	0.00
20		Hexachlorobenzene	1.000	1.167	-16.7	133	0.00
21		Pentachlorophenol	1.000	1.123	-12.3	125	0.00
22		Phenanthrene	1.000	0.972	2.8	98	0.00
23		Anthracene	1.000	1.224	-22.4	140	0.00
24		Fluoranthene	1.000	1.014	-1.4	117	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	116	0.00
26		Pyrene	1.000	0.967	3.3	118	0.00
27	S	Terphenyl-d14	1.000	0.975	2.5	117	0.00
28		Benzo(a)anthracene	1.000	0.928	7.2	114	0.00
29		Chrysene	1.000	0.991	0.9	116	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.849	15.1	110	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.923	7.7	112	0.00
32	I	Perylene-d12	5.000	5.000	0.0	114	0.01
33		Benzo(b)fluoranthene	1.000	0.960	4.0	114	0.00
34		Benzo(k)fluoranthene	1.000	1.000	0.0	115	0.00
35	C	Benzo(a)pyrene	1.000	0.961	3.9	113	0.00
36		Dibenzo(a,h)anthracene	1.000	0.945	5.5	112	0.00
37		Benzo(g,h,i)perylene	1.000	0.942	5.8	112	0.00

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