Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM060615\
Data File : BM001579.D
Acq On : 06 Jun 2015 05:42
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 08 05:43:25 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 08 05:27:25 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	103	0.00
2		1,4-Dioxane	1.000	0.922	7.8	102	0.00
3		n-Nitrosodimethylamine	1.000	0.972	2.8	103	0.00
4	S	2-Fluorophenol	1.000	0.952	4.8	102	0.00
5	S	Phenol-d6	1.000	0.959	4.1	103	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	104	0.00
7	S	Nitrobenzene-d5	1.000	0.943	5.7	102	0.00
8		Nitrobenzene	1.000	0.949	5.1	103	0.00
9		Naphthalene	1.000	0.984	1.6	105	0.00
10		Hexachlorobutadiene	1.000	0.975	2.5	105	0.00
11		2-Methylnaphthalene	1.000	0.979	2.1	105	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	107	0.00
13	S	2,4,6-Tribromophenol	1.000	0.919	8.1	112	0.00
14	S	2-Fluorobiphenyl	1.000	0.946	5.4	105	0.00
15		Acenaphthylene	1.000	0.950	5.0	105	0.00
16		Acenaphthene	1.000	0.952	4.8	106	0.00
17		Fluorene	1.000	0.966	3.4	111	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	83	0.00
19		4-Bromophenyl-phenylether	1.000	1.122	-12.2	115	0.00
20		Hexachlorobenzene	1.000	1.377	-37.7#	127	0.00
21		Pentachlorophenol	1.000	0.942	5.8	105	0.00
22		Phenanthrene	1.000	0.984	1.6	87	0.00
23		Anthracene	1.000	1.196	-19.6	116	0.00
24		Fluoranthene	1.000	1.229	-22.9	113	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
26		Pyrene	1.000	0.977	2.3	113	0.00
27	S	Terphenyl-d14	1.000	0.989	1.1	114	0.00
28		Benzo(a)anthracene	1.000	0.945	5.5	114	0.00
29		Chrysene	1.000	0.960	4.0	114	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.946	5.4	121	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.777	22.3	96	0.00
32	I	Perylene-d12	5.000	5.000	0.0	104	0.00
33		Benzo(b)fluoranthene	1.000	0.930	7.0	106	0.00
34		Benzo(k)fluoranthene	1.000	1.001	-0.1	103	0.00
35	C	Benzo(a)pyrene	1.000	0.927	7.3	102	0.00
36		Dibenzo(a,h)anthracene	1.000	0.846	15.4	96	0.00
37		Benzo(g,h,i)perylene	1.000	0.847	15.3	96	0.00

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