Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM050115\
Data File : BM001153.D
Acq On : 02 May 2015 04:31
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 04 17:09:06 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM050115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 04 16:48:07 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	98	0.00
2		1,4-Dioxane	1.000	1.071	-7.1	94	0.02
3		n-Nitrosodimethylamine	1.000	0.919	8.1	85	0.00
4	S	2-Fluorophenol	1.000	1.009	-0.9	97	0.00
5	S	Phenol-d6	1.000	0.996	0.4	97	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
7	S	Nitrobenzene-d5	1.000	0.988	1.2	97	0.00
8		Nitrobenzene	1.000	0.981	1.9	96	0.00
9		Naphthalene	1.000	1.010	-1.0	96	0.00
10		Hexachlorobutadiene	1.000	1.011	-1.1	99	0.00
11		2-Methylnaphthalene	1.000	1.024	-2.4	100	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	1.000	1.267	-26.7#	157	0.00
14	S	2-Fluorobiphenyl	1.000	1.032	-3.2	101	0.00
15		Acenaphthylene	1.000	1.026	-2.6	101	-0.02
16		Acenaphthene	1.000	1.025	-2.5	103	0.00
17		Fluorene	1.000	1.039	-3.9	100	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
19		Hexachlorobenzene	1.000	0.990	1.0	102	0.00
20		Pentachlorophenol	1.000	0.858	14.2	95	0.00
21		Phenanthrene	1.000	0.986	1.4	101	0.00
22		Anthracene	1.000	1.004	-0.4	100	0.00
23		Fluoranthene	1.000	1.001	-0.1	102	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
25		Pyrene	1.000	0.995	0.5	101	0.00
26	S	Terphenyl-d14	1.000	1.015	-1.5	101	0.00
27		Benzo(a)anthracene	1.000	0.999	0.1	102	0.00
28		Chrysene	1.000	1.001	-0.1	98	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	1.023	-2.3	100	0.00
30	I	Perylene-d12	5.000	5.000	0.0	100	-0.01
31		Benzo(b)fluoranthene	1.000	0.989	1.1	92	-0.01
32		Benzo(k)fluoranthene	1.000	1.046	-4.6	112	-0.01
33	C	Benzo(a)pyrene	1.000	1.013	-1.3	101	0.00
34		Dibenzo(a,h)anthracene	1.000	1.020	-2.0	101	-0.01
35		Benzo(g,h,i)perylene	1.000	1.010	-1.0	99	0.00

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