Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM051615\
Data File : BM001294.D
Acq On : 14 May 2015 16:06
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 15 05:26:32 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 05:24:15 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	97	0.00
2		1,4-Dioxane	1.000	0.832	16.8	88	0.00
3		n-Nitrosodimethylamine	1.000	0.991	0.9	110	0.00
4	S	2-Fluorophenol	1.000	0.944	5.6	97	0.00
5	S	Phenol-d6	1.000	0.941	5.9	99	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	96	0.00
7	S	Nitrobenzene-d5	1.000	0.900	10.0	98	0.00
8		Nitrobenzene	1.000	0.875	12.5	97	0.00
9		Naphthalene	1.000	0.978	2.2	97	0.00
10		Hexachlorobutadiene	1.000	0.964	3.6	98	0.00
11		2-Methylnaphthalene	1.000	0.948	5.2	97	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.804	19.6	95	0.00
14	S	2-Fluorobiphenyl	1.000	0.949	5.1	95	0.00
15		Acenaphthylene	1.000	0.943	5.7	97	0.00
16		Acenaphthene	1.000	0.945	5.5	97	0.00
17		Fluorene	1.000	0.852	14.8	92	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.00
19		Hexachlorobenzene	1.000	0.857	14.3	100	0.00
20		Pentachlorophenol	1.000	0.847	15.3	97	0.00
21		Phenanthrene	1.000	0.855	14.5	100	0.00
22		Anthracene	1.000	0.832	16.8	98	0.00
23		Fluoranthene	1.000	0.814	18.6	97	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	97	0.00
25		Pyrene	1.000	0.920	8.0	98	0.00
26	S	Terphenyl-d14	1.000	0.900	10.0	93	-0.01
27		Benzo(a)anthracene	1.000	0.905	9.5	99	0.00
28		Chrysene	1.000	0.950	5.0	98	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	1.020	-2.0	107	0.00
30	I	Perylene-d12	5.000	5.000	0.0	101	-0.01
31		Benzo(b)fluoranthene	1.000	0.928	7.2	102	0.00
32		Benzo(k)fluoranthene	1.000	0.928	7.2	100	0.00
33	C	Benzo(a)pyrene	1.000	0.938	6.2	104	0.00
34		Dibenzo(a,h)anthracene	1.000	0.993	0.7	106	0.00
35		Benzo(g,h,i)perylene	1.000	0.991	0.9	107	-0.01

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