Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051915\
Data File : BM001344.D
Acq On : 19 May 2015 12:26
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 20 04:10:26 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 20 03:51:55 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	96	-0.02
2		1,4-Dioxane	1.000	0.987	1.3	95	0.00
3		n-Nitrosodimethylamine	1.000	1.010	-1.0	111	0.00
4	S	2-Fluorophenol	1.000	0.988	1.2	100	-0.02
5	S	Phenol-d6	1.000	1.007	-0.7	105	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
7	S	Nitrobenzene-d5	1.000	0.939	6.1	106	0.00
8		Nitrobenzene	1.000	0.921	7.9	107	0.00
9		Naphthalene	1.000	1.013	-1.3	104	0.00
10		Hexachlorobutadiene	1.000	1.011	-1.1	106	0.00
11		2-Methylnaphthalene	1.000	1.035	-3.5	109	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	1.000	0.719	28.1#	90	0.00
14	S	2-Fluorobiphenyl	1.000	1.023	-2.3	111	0.00
15		Acenaphthylene	1.000	0.986	1.4	110	0.00
16		Acenaphthene	1.000	0.988	1.2	110	0.00
17		Fluorene	1.000	0.927	7.3	108	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	129	0.00
19		Hexachlorobenzene	1.000	0.728	27.2#	102	0.00
20		Pentachlorophenol	1.000	0.935	6.5	140	0.00
21		Phenanthrene	1.000	1.116	-11.6	157	0.00
22		Anthracene	1.000	1.028	-2.8	145	0.00
23		Fluoranthene	1.000	0.687	31.3#	98	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	88	-0.01
25		Pyrene	1.000	1.019	-1.9	99	0.00
26	S	Terphenyl-d14	1.000	1.045	-4.5	99	-0.01
27		Benzo(a)anthracene	1.000	0.942	5.8	94	0.00
28		Chrysene	1.000	0.985	1.5	92	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	0.965	3.5	92	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	88	-0.01
31		Benzo(b)fluoranthene	1.000	1.048	-4.8	100	0.00
32		Benzo(k)fluoranthene	1.000	0.886	11.4	83	-0.01
33	C	Benzo(a)pyrene	1.000	0.944	5.6	91	0.00
34		Dibenzo(a,h)anthracene	1.000	0.978	2.2	91	-0.01
35		Benzo(g,h,i)perylene	1.000	0.989	1.1	93	-0.02

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