Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM051815\
Data File : BM001340.D
Acq On : 18 May 2015 20:59
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 19 02:32:58 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 19 02:11:49 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	101	0.00
2		1,4-Dioxane	1.000	0.403	59.7#	66	-0.01
3		n-Nitrosodimethylamine	1.000	0.927	7.3	107	0.00
4	S	2-Fluorophenol	1.000	0.937	6.3	100	-0.02
5	S	Phenol-d6	1.000	0.969	3.1	106	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	102	0.00
7	S	Nitrobenzene-d5	1.000	0.913	8.7	106	0.00
8		Nitrobenzene	1.000	0.880	12.0	105	0.00
9		Naphthalene	1.000	0.964	3.6	102	0.00
10		Hexachlorobutadiene	1.000	0.954	4.6	104	0.00
11		2-Methylnaphthalene	1.000	0.956	4.4	104	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
13	S	2,4,6-Tribromophenol	1.000	0.700	30.0#	86	0.00
14	S	2-Fluorobiphenyl	1.000	0.962	3.8	103	0.00
15		Acenaphthylene	1.000	0.947	5.3	103	0.00
16		Acenaphthene	1.000	0.947	5.3	103	0.00
17		Fluorene	1.000	0.847	15.3	97	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	130	0.00
19		Hexachlorobenzene	1.000	0.714	28.6#	100	0.00
20		Pentachlorophenol	1.000	0.909	9.1	134	0.00
21		Phenanthrene	1.000	0.959	4.1	136	0.00
22		Anthracene	1.000	0.927	7.3	132	0.00
23		Fluoranthene	1.000	0.704	29.6#	101	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	-0.01
25		Pyrene	1.000	0.928	7.2	103	0.00
26	S	Terphenyl-d14	1.000	0.955	4.5	103	-0.01
27		Benzo(a)anthracene	1.000	0.923	7.7	105	-0.01
28		Chrysene	1.000	0.947	5.3	102	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	0.941	5.9	102	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	101	-0.01
31		Benzo(b)fluoranthene	1.000	0.906	9.4	100	-0.01
32		Benzo(k)fluoranthene	1.000	0.976	2.4	105	-0.01
33	C	Benzo(a)pyrene	1.000	0.930	7.0	103	-0.01
34		Dibenzo(a,h)anthracene	1.000	0.949	5.1	102	-0.01
35		Benzo(g,h,i)perylene	1.000	0.942	5.8	102	-0.02

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