Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051815\
Data File : BM001327.D
Acq On : 15 May 2015 14:26
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 16 01:49:33 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	94	-0.02
2		1,4-Dioxane	1.000	0.925	7.5	90	0.00
3		n-Nitrosodimethylamine	1.000	1.004	-0.4	108	0.00
4	S	2-Fluorophenol	1.000	0.962	3.8	95	-0.02
5	S	Phenol-d6	1.000	0.948	5.2	96	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	93	0.00
7	S	Nitrobenzene-d5	1.000	0.876	12.4	92	0.00
8		Nitrobenzene	1.000	0.866	13.4	93	0.00
9		Naphthalene	1.000	0.969	3.1	93	0.00
10		Hexachlorobutadiene	1.000	0.970	3.0	96	0.00
11		2-Methylnaphthalene	1.000	0.928	7.2	92	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	90	0.00
13	S	2,4,6-Tribromophenol	1.000	0.685	31.5#	74	-0.03
14	S	2-Fluorobiphenyl	1.000	0.967	3.3	91	0.00
15		Acenaphthylene	1.000	0.949	5.1	92	0.00
16		Acenaphthene	1.000	0.952	4.8	92	0.00
17		Fluorene	1.000	0.963	3.7	97	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	121	0.00
19		Hexachlorobenzene	1.000	0.664	33.6#	87	0.00
20		Pentachlorophenol	1.000	0.914	8.6	126	0.00
21		Phenanthrene	1.000	1.119	-11.9	148	0.00
22		Anthracene	1.000	0.993	0.7	131	0.00
23		Fluoranthene	1.000	0.701	29.9#	94	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	96	-0.01
25		Pyrene	1.000	0.923	7.7	97	0.00
26	S	Terphenyl-d14	1.000	0.918	8.2	94	-0.01
27		Benzo(a)anthracene	1.000	0.926	7.4	100	-0.01
28		Chrysene	1.000	0.947	5.3	96	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	1.026	-2.6	106	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	100	-0.01
31		Benzo(b)fluoranthene	1.000	0.901	9.9	99	-0.01
32		Benzo(k)fluoranthene	1.000	0.979	2.1	105	-0.01
33	C	Benzo(a)pyrene	1.000	0.944	5.6	104	-0.01
34		Dibenzo(a,h)anthracene	1.000	0.992	0.8	106	-0.01
35		Benzo(g,h,i)perylene	1.000	0.989	1.1	106	-0.02

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