Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM051615\
Data File : BM001309.D
Acq On : 15 May 2015 01:32
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 15 05:34:12 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 15 05:14:53 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	0.668	33.2#	80	0.00
3		n-Nitrosodimethylamine	1.000	0.984	1.6	111	0.00
4	S	2-Fluorophenol	1.000	0.946	5.4	98	-0.02
5	S	Phenol-d6	1.000	0.948	5.2	101	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
7	S	Nitrobenzene-d5	1.000	0.874	12.6	98	0.00
8		Nitrobenzene	1.000	0.848	15.2	97	0.00
9		Naphthalene	1.000	0.963	3.7	99	0.00
10		Hexachlorobutadiene	1.000	0.934	6.6	99	0.00
11		2-Methylnaphthalene	1.000	0.936	6.4	99	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
13	S	2,4,6-Tribromophenol	1.000	0.720	28.0#	85	0.00
14	S	2-Fluorobiphenyl	1.000	0.951	4.9	97	0.00
15		Acenaphthylene	1.000	0.933	6.7	97	0.00
16		Acenaphthene	1.000	0.938	6.2	98	0.00
17		Fluorene	1.000	0.916	8.4	100	-0.03
18	I	Phenanthrene-d10	5.000	5.000	0.0	132	0.00
19		Hexachlorobenzene	1.000	0.659	34.1#	94	0.00
20		Pentachlorophenol	1.000	0.916	8.4	138	0.00
21		Phenanthrene	1.000	0.977	2.3	140	0.00
22		Anthracene	1.000	0.937	6.3	135	0.00
23		Fluoranthene	1.000	0.680	32.0#	99	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	100	-0.01
25		Pyrene	1.000	0.917	8.3	100	0.00
26	S	Terphenyl-d14	1.000	0.904	9.6	97	-0.01
27		Benzo(a)anthracene	1.000	0.934	6.6	105	-0.01
28		Chrysene	1.000	0.934	6.6	99	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	1.034	-3.4	111	-0.01
30	I	Perylene-d12	5.000	5.000	0.0	104	-0.02
31		Benzo(b)fluoranthene	1.000	0.930	7.0	106	-0.01
32		Benzo(k)fluoranthene	1.000	0.932	6.8	104	-0.01
33	C	Benzo(a)pyrene	1.000	0.938	6.2	108	-0.01
34		Dibenzo(a,h)anthracene	1.000	0.996	0.4	111	-0.02
35		Benzo(g,h,i)perylene	1.000	0.987	1.3	111	-0.02

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