Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM050115\
Data File : BM001170.D
Acq On : 02 May 2015 15:50
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 04 17:35:54 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM050115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 04 16:48:07 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	0.697	30.3#	73	0.00
3		n-Nitrosodimethylamine	1.000	0.983	1.7	94	0.00
4	S	2-Fluorophenol	1.000	1.015	-1.5	101	-0.01
5	S	Phenol-d6	1.000	1.014	-1.4	101	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
7	S	Nitrobenzene-d5	1.000	0.985	1.5	99	0.00
8		Nitrobenzene	1.000	0.933	6.7	93	0.00
9		Naphthalene	1.000	1.009	-0.9	99	0.00
10		Hexachlorobutadiene	1.000	1.017	-1.7	102	0.00
11		2-Methylnaphthalene	1.000	1.043	-4.3	104	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	1.000	1.280	-28.0#	164	0.00
14	S	2-Fluorobiphenyl	1.000	1.017	-1.7	103	0.00
15		Acenaphthylene	1.000	1.028	-2.8	104	-0.02
16		Acenaphthene	1.000	1.016	-1.6	106	0.00
17		Fluorene	1.000	1.055	-5.5	104	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
19		Hexachlorobenzene	1.000	1.016	-1.6	107	0.00
20		Pentachlorophenol	1.000	0.879	12.1	101	0.00
21		Phenanthrene	1.000	1.039	-3.9	108	0.00
22		Anthracene	1.000	0.985	1.5	100	0.00
23		Fluoranthene	1.000	1.006	-0.6	105	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
25		Pyrene	1.000	1.034	-3.4	104	0.00
26	S	Terphenyl-d14	1.000	1.038	-3.8	103	0.00
27		Benzo(a)anthracene	1.000	1.001	-0.1	102	0.00
28		Chrysene	1.000	1.006	-0.6	99	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	0.890	11.0	87	0.00
30	I	Perylene-d12	5.000	5.000	0.0	93	-0.01
31		Benzo(b)fluoranthene	1.000	0.998	0.2	86	-0.01
32		Benzo(k)fluoranthene	1.000	1.071	-7.1	106	-0.01
33	C	Benzo(a)pyrene	1.000	1.020	-2.0	95	0.00
34		Dibenzo(a,h)anthracene	1.000	0.952	4.8	87	-0.01
35		Benzo(g,h,i)perylene	1.000	0.937	6.3	86	0.00

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