Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050415\
Data File : BM001186.D
Acq On : 04 May 2015 10:52
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 05 02:43:47 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	128	0.00
2		1,4-Dioxane	1.000	0.883	11.7	108	0.00
3		n-Nitrosodimethylamine	1.000	1.016	-1.6	124	0.00
4	S	2-Fluorophenol	1.000	1.043	-4.3	132	-0.01
5	S	Phenol-d6	1.000	1.079	-7.9	137	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	132	0.00
7	S	Nitrobenzene-d5	1.000	1.019	-1.9	135	0.00
8		Nitrobenzene	1.000	0.959	4.1	127	0.00
9		Naphthalene	1.000	1.003	-0.3	130	0.00
10		Hexachlorobutadiene	1.000	0.998	0.2	132	0.00
11		2-Methylnaphthalene	1.000	1.035	-3.5	137	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	136	0.00
13	S	2,4,6-Tribromophenol	1.000	1.006	-0.6	164	0.00
14	S	2-Fluorobiphenyl	1.000	1.024	-2.4	137	-0.01
15		Acenaphthylene	1.000	1.067	-6.7	143	-0.02
16		Acenaphthene	1.000	1.058	-5.8	146	0.00
17		Fluorene	1.000	1.163	-16.3	152	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	144	0.00
19		Hexachlorobenzene	1.000	0.977	2.3	143	0.00
20		Pentachlorophenol	1.000	0.942	5.8	157	0.00
21		Phenanthrene	1.000	1.016	-1.6	147	0.00
22		Anthracene	1.000	1.010	-1.0	142	-0.02
23		Fluoranthene	1.000	1.021	-2.1	147	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	137	0.00
25		Pyrene	1.000	1.072	-7.2	147	0.00
26	S	Terphenyl-d14	1.000	1.059	-5.9	142	0.00
27		Benzo(a)anthracene	1.000	1.051	-5.1	145	0.00
28		Chrysene	1.000	1.014	-1.4	134	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	0.880	12.0	116	0.00
30	I	Perylene-d12	5.000	5.000	0.0	124	-0.01
31		Benzo(b)fluoranthene	1.000	1.004	-0.4	115	0.00
32		Benzo(k)fluoranthene	1.000	1.111	-11.1	146	-0.01
33	C	Benzo(a)pyrene	1.000	1.055	-5.5	130	0.00
34		Dibenzo(a,h)anthracene	1.000	0.949	5.1	115	-0.01
35		Benzo(g,h,i)perylene	1.000	0.933	6.7	113	0.00

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