Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM052015\
Data File : BM001363.D
Acq On : 21 May 2015 16:20
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 21 17:05:38 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM052015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 21 11:08:14 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	99	0.00
2		1,4-Dioxane	1.000	1.359	-35.9#	113	0.00
3		n-Nitrosodimethylamine	1.000	1.097	-9.7	102	0.00
4	S	2-Fluorophenol	1.000	1.052	-5.2	105	0.00
5	S	Phenol-d6	1.000	1.080	-8.0	108	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
7	S	Nitrobenzene-d5	1.000	1.058	-5.8	109	0.00
8		Nitrobenzene	1.000	0.996	0.4	111	0.00
9		Naphthalene	1.000	1.046	-4.6	104	0.00
10		Hexachlorobutadiene	1.000	1.045	-4.5	105	0.00
11		2-Methylnaphthalene	1.000	1.057	-5.7	106	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	103	0.00
13	S	2,4,6-Tribromophenol	1.000	0.950	5.0	101	0.00
14	S	2-Fluorobiphenyl	1.000	1.042	-4.2	107	0.00
15		Acenaphthylene	1.000	1.046	-4.6	111	0.00
16		Acenaphthene	1.000	1.045	-4.5	109	0.00
17		Fluorene	1.000	1.040	-4.0	109	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
19		4-Bromophenyl-phenylether	1.000	0.900	10.0	105	0.00
20		Hexachlorobenzene	1.000	0.966	3.4	109	0.00
21		Pentachlorophenol	1.000	0.959	4.1	131	0.00
22		Phenanthrene	1.000	0.999	0.1	111	0.00
23		Anthracene	1.000	0.994	0.6	104	0.00
24		Fluoranthene	1.000	0.991	0.9	110	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	106	0.00
26		Pyrene	1.000	1.032	-3.2	111	0.00
27	S	Terphenyl-d14	1.000	1.030	-3.0	112	0.00
28		Benzo(a)anthracene	1.000	1.006	-0.6	113	0.00
29		Chrysene	1.000	1.024	-2.4	112	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.057	-5.7	134	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.015	-1.5	116	0.00
32	I	Perylene-d12	5.000	5.000	0.0	109	0.00
33		Benzo(b)fluoranthene	1.000	0.989	1.1	110	0.00
34		Benzo(k)fluoranthene	1.000	1.074	-7.4	118	0.00
35	C	Benzo(a)pyrene	1.000	1.037	-3.7	118	0.00
36		Dibenzo(a,h)anthracene	1.000	1.021	-2.1	116	0.01
37		Benzo(g,h,i)perylene	1.000	1.019	-1.9	117	0.00

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