Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061915\
Data File : BM001754.D
Acq On : 19 Jun 2015 10:15
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 19 12:24:30 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 19 05:05:14 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	110	0.00
2		1,4-Dioxane	1.000	1.009	-0.9	107	0.00
3		n-Nitrosodimethylamine	1.000	1.033	-3.3	113	-0.02
4	S	2-Fluorophenol	1.000	1.005	-0.5	110	0.00
5	S	Phenol-d6	1.000	0.994	0.6	111	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	110	-0.02
7	S	Nitrobenzene-d5	1.000	0.990	1.0	111	0.00
8		Nitrobenzene	1.000	1.000	0.0	111	0.00
9		Naphthalene	1.000	0.990	1.0	109	0.00
10		Hexachlorobutadiene	1.000	0.991	0.9	112	0.00
11		2-Methylnaphthalene	1.000	0.989	1.1	110	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	109	-0.01
13	S	2,4,6-Tribromophenol	1.000	1.092	-9.2	122	0.00
14	S	2-Fluorobiphenyl	1.000	1.003	-0.3	109	0.00
15		Acenaphthylene	1.000	0.992	0.8	110	0.00
16		Acenaphthene	1.000	0.994	0.6	110	0.00
17		Fluorene	1.000	1.017	-1.7	99	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
19		4-Bromophenyl-phenylether	1.000	1.153	-15.3	124	0.00
20		Hexachlorobenzene	1.000	0.959	4.1	100	0.00
21		Pentachlorophenol	1.000	1.036	-3.6	109	0.00
22		Phenanthrene	1.000	0.986	1.4	101	0.00
23		Anthracene	1.000	1.025	-2.5	100	0.00
24		Fluoranthene	1.000	1.088	-8.8	110	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	109	0.00
26		Pyrene	1.000	0.965	3.5	111	-0.01
27	S	Terphenyl-d14	1.000	0.990	1.0	112	0.00
28		Benzo(a)anthracene	1.000	0.974	2.6	109	0.00
29		Chrysene	1.000	0.976	2.4	108	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.948	5.2	113	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.981	1.9	107	0.00
32	I	Perylene-d12	5.000	5.000	0.0	109	0.00
33		Benzo(b)fluoranthene	1.000	1.009	-0.9	111	-0.01
34		Benzo(k)fluoranthene	1.000	0.962	3.8	107	0.00
35	C	Benzo(a)pyrene	1.000	0.986	1.4	109	0.00
36		Dibenzo(a,h)anthracene	1.000	0.978	2.2	106	0.00
37		Benzo(g,h,i)perylene	1.000	0.978	2.2	107	0.00

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