Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052215\
Data File : BM001411.D
Acq On : 23 May 2015 09:23
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 25 08:28:05 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 23 04:45:39 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	125	0.00
2		1,4-Dioxane	1.000	1.217	-21.7	154	-0.02
3		n-Nitrosodimethylamine	1.000	1.028	-2.8	126	-0.02
4	S	2-Fluorophenol	1.000	1.028	-2.8	124	0.00
5	S	Phenol-d6	1.000	1.051	-5.1	128	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	128	-0.02
7	S	Nitrobenzene-d5	1.000	1.002	-0.2	128	0.00
8		Nitrobenzene	1.000	1.059	-5.9	136	0.00
9		Naphthalene	1.000	1.046	-4.6	130	0.00
10		Hexachlorobutadiene	1.000	1.041	-4.1	128	0.00
11		2-Methylnaphthalene	1.000	1.063	-6.3	132	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	133	0.00
13	S	2,4,6-Tribromophenol	1.000	0.719	28.1#	94	0.00
14	S	2-Fluorobiphenyl	1.000	1.031	-3.1	132	0.00
15		Acenaphthylene	1.000	1.063	-6.3	138	0.00
16		Acenaphthene	1.000	1.031	-3.1	134	0.00
17		Fluorene	1.000	0.830	17.0	113	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	135	0.00
19		4-Bromophenyl-phenylether	1.000	1.032	-3.2	114	0.00
20		Hexachlorobenzene	1.000	1.249	-24.9	166	0.00
21		Pentachlorophenol	1.000	1.159	-15.9	129	-0.03
22		Phenanthrene	1.000	1.244	-24.4	163	0.00
23		Anthracene	1.000	0.856	14.4	104	0.00
24		Fluoranthene	1.000	0.918	8.2	127	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	114	0.00
26		Pyrene	1.000	1.114	-11.4	125	0.00
27	S	Terphenyl-d14	1.000	1.068	-6.8	119	-0.01
28		Benzo(a)anthracene	1.000	1.023	-2.3	118	0.00
29		Chrysene	1.000	0.986	1.4	110	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.886	11.4	112	-0.02
31		Indeno(1,2,3-cd)pyrene	1.000	0.856	14.4	95	-0.01
32	I	Perylene-d12	5.000	5.000	0.0	102	0.00
33		Benzo(b)fluoranthene	1.000	1.052	-5.2	109	0.00
34		Benzo(k)fluoranthene	1.000	1.034	-3.4	101	0.00
35	C	Benzo(a)pyrene	1.000	1.010	-1.0	103	0.00
36		Dibenzo(a,h)anthracene	1.000	0.928	7.2	93	-0.01
37		Benzo(g,h,i)perylene	1.000	0.970	3.0	97	-0.01

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