Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030516\
Data File : BM004571.D
Acq On : 06 Mar 2016 00:07
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Mar 07 08:04:39 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM030516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 07 07:54:56 2016
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	0.946	5.4	109	0.00
3		n-Nitrosodimethylamine	1.000	0.866	13.4	106	0.00
4	S	2-Fluorophenol	1.000	0.829	17.1	101	0.02
5	S	Phenol-d6	1.000	0.841	15.9	102	0.00
6		bis(2-Chloroethyl)ether	1.000	0.859	14.1	102	0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
8	S	Nitrobenzene-d5	1.000	0.853	14.7	101	0.00
9		Nitrobenzene	1.000	0.835	16.5	101	0.00
10		Naphthalene	1.000	0.997	0.3	118	0.01
11		Hexachlorobutadiene	1.000	0.917	8.3	108	0.00
12		2-Methylnaphthalene	1.000	0.884	11.6	105	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	100	-0.02
14	S	2,4,6-Tribromophenol	1.000	0.864	13.6	96	0.00
15	S	2-Fluorobiphenyl	1.000	0.991	0.9	105	0.00
16		Acenaphthylene	1.000	0.942	5.8	101	0.00
17		Acenaphthene	1.000	1.046	-4.6	112	-0.02
18		Fluorene	1.000	0.923	7.7	99	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	90	0.00
20		Hexachlorobenzene	1.000	1.051	-5.1	98	-0.02
21		Pentachlorophenol	1.000	0.611	38.9#	68	0.00
22		Phenanthrene	1.000	1.171	-17.1	111	-0.02
23		Anthracene	1.000	1.072	-7.2	112	-0.02
24		Fluoranthene	1.000	1.069	-6.9	103	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	106	-0.01
26		Benzidine	1.000	1.105	-10.5	161	0.00
27		Pyrene	1.000	0.893	10.7	103	0.00
28	S	Terphenyl-d14	1.000	0.882	11.8	104	0.00
29		Benzo(a)anthracene	1.000	0.887	11.3	104	0.00
30		3,3'-Dichlorobenzidine	1.000	0.884	11.6	99	0.01
31		Chrysene	1.000	0.918	8.2	103	-0.01
32		Bis(2-ethylhexyl)phthalate	1.000	0.862	13.8	105	-0.01
33		Indeno(1,2,3-cd)pyrene	1.000	0.899	10.1	104	0.00
34	I	Perylene-d12	5.000	5.000	0.0	107	0.00
35	C	Benzo(a)pyrene	1.000	0.905	9.5	106	0.00
36		Dibenzo(a,h)anthracene	1.000	0.857	14.3	105	0.00
37		Benzo(g,h,i)perylene	1.000	0.867	13.3	105	0.00

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