Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061315\
Data File : BM001693.D
Acq On : 13 Jun 2015 05:24
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jun 13 06:49:56 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jun 13 02:11:48 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	95	0.00
2		1,4-Dioxane	1.000	1.094	-9.4	96	0.00
3		n-Nitrosodimethylamine	1.000	0.993	0.7	96	0.00
4	S	2-Fluorophenol	1.000	1.005	-0.5	97	0.00
5	S	Phenol-d6	1.000	1.006	-0.6	96	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
7	S	Nitrobenzene-d5	1.000	0.987	1.3	95	0.00
8		Nitrobenzene	1.000	0.981	1.9	96	0.00
9		Naphthalene	1.000	0.991	0.9	95	0.00
10		Hexachlorobutadiene	1.000	0.985	1.5	94	0.00
11		2-Methylnaphthalene	1.000	0.991	0.9	95	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	1.067	-6.7	106	0.00
14	S	2-Fluorobiphenyl	1.000	1.008	-0.8	96	0.00
15		Acenaphthylene	1.000	0.995	0.5	95	0.00
16		Acenaphthene	1.000	0.987	1.3	94	0.00
17		Fluorene	1.000	0.928	7.2	93	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	93	0.00
19		4-Bromophenyl-phenylether	1.000	1.211	-21.1	111	0.00
20		Hexachlorobenzene	1.000	0.900	10.0	94	0.00
21		Pentachlorophenol	1.000	0.941	5.9	91	0.00
22		Phenanthrene	1.000	1.164	-16.4	102	0.00
23		Anthracene	1.000	0.903	9.7	91	0.00
24		Fluoranthene	1.000	1.011	-1.1	99	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
26		Pyrene	1.000	1.001	-0.1	101	0.00
27	S	Terphenyl-d14	1.000	0.977	2.3	97	0.00
28		Benzo(a)anthracene	1.000	1.000	0.0	102	0.00
29		Chrysene	1.000	1.017	-1.7	100	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.035	-3.5	107	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.004	-0.4	100	0.00
32	I	Perylene-d12	5.000	5.000	0.0	100	0.00
33		Benzo(b)fluoranthene	1.000	0.999	0.1	104	0.00
34		Benzo(k)fluoranthene	1.000	1.025	-2.5	99	-0.01
35	C	Benzo(a)pyrene	1.000	1.008	-0.8	101	0.00
36		Dibenzo(a,h)anthracene	1.000	1.009	-0.9	100	0.00
37		Benzo(g,h,i)perylene	1.000	0.996	0.4	100	0.00

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