Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM051315\
Data File : BM001274.D
Acq On : 13 May 2015 05:04
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 13 06:46:32 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM051315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 13 05:24:15 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	94	0.00
2		1,4-Dioxane	1.000	1.043	-4.3	97	0.00
3		n-Nitrosodimethylamine	1.000	0.884	11.6	96	0.00
4	S	2-Fluorophenol	1.000	0.948	5.2	95	0.00
5	S	Phenol-d6	1.000	0.936	6.4	96	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	1.000	0.884	11.6	94	0.00
8		Nitrobenzene	1.000	0.862	13.8	94	0.00
9		Naphthalene	1.000	0.973	2.7	95	0.00
10		Hexachlorobutadiene	1.000	0.950	5.0	95	0.00
11		2-Methylnaphthalene	1.000	0.946	5.4	95	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
13	S	2,4,6-Tribromophenol	1.000	0.750	25.0	86	0.00
14	S	2-Fluorobiphenyl	1.000	0.958	4.2	94	0.00
15		Acenaphthylene	1.000	0.932	6.8	93	0.00
16		Acenaphthene	1.000	0.944	5.6	94	0.00
17		Fluorene	1.000	0.832	16.8	87	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
19		Hexachlorobenzene	1.000	0.791	20.9	96	0.00
20		Pentachlorophenol	1.000	0.914	8.6	117	0.00
21		Phenanthrene	1.000	0.818	18.2	100	0.00
22		Anthracene	1.000	0.823	17.7	101	0.00
23		Fluoranthene	1.000	0.752	24.8	94	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	97	-0.01
25		Pyrene	1.000	0.910	9.0	97	0.00
26	S	Terphenyl-d14	1.000	0.906	9.4	94	-0.01
27		Benzo(a)anthracene	1.000	0.889	11.1	98	0.00
28		Chrysene	1.000	0.947	5.3	98	-0.01
29		Indeno(1,2,3-cd)pyrene	1.000	0.992	0.8	103	0.00
30	I	Perylene-d12	5.000	5.000	0.0	100	-0.01
31		Benzo(b)fluoranthene	1.000	0.969	3.1	106	0.00
32		Benzo(k)fluoranthene	1.000	0.857	14.3	91	0.00
33	C	Benzo(a)pyrene	1.000	0.909	9.1	100	0.00
34		Dibenzo(a,h)anthracene	1.000	0.968	3.2	103	-0.01
35		Benzo(g,h,i)perylene	1.000	0.977	2.3	104	-0.01

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