Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052215\
Data File : BM001395.D
Acq On : 22 May 2015 23:16
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 23 05:14:18 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	97	0.00
2		1,4-Dioxane	1.000	1.098	-9.8	110	0.00
3		n-Nitrosodimethylamine	1.000	1.110	-11.0	106	-0.02
4	S	2-Fluorophenol	1.000	1.027	-2.7	96	0.00
5	S	Phenol-d6	1.000	1.025	-2.5	97	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
7	S	Nitrobenzene-d5	1.000	0.983	1.7	96	0.00
8		Nitrobenzene	1.000	0.977	2.3	97	0.00
9		Naphthalene	1.000	1.029	-2.9	98	0.00
10		Hexachlorobutadiene	1.000	1.042	-4.2	98	0.00
11		2-Methylnaphthalene	1.000	1.039	-3.9	99	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	100	0.00
13	S	2,4,6-Tribromophenol	1.000	0.791	20.9	78	0.00
14	S	2-Fluorobiphenyl	1.000	1.029	-2.9	99	0.00
15		Acenaphthylene	1.000	1.030	-3.0	100	0.00
16		Acenaphthene	1.000	1.024	-2.4	99	0.00
17		Fluorene	1.000	0.922	7.8	94	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
19		4-Bromophenyl-phenylether	1.000	1.035	-3.5	86	0.00
20		Hexachlorobenzene	1.000	1.221	-22.1	122	0.00
21		Pentachlorophenol	1.000	1.179	-17.9	100	0.00
22		Phenanthrene	1.000	1.200	-20.0	118	0.00
23		Anthracene	1.000	0.945	5.5	87	0.00
24		Fluoranthene	1.000	0.987	1.3	103	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	104	0.00
26		Pyrene	1.000	1.005	-0.5	103	0.00
27	S	Terphenyl-d14	1.000	0.993	0.7	101	-0.01
28		Benzo(a)anthracene	1.000	1.049	-4.9	110	0.00
29		Chrysene	1.000	0.945	5.5	96	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.851	14.9	98	-0.02
31		Indeno(1,2,3-cd)pyrene	1.000	0.990	1.0	100	-0.01
32	I	Perylene-d12	5.000	5.000	0.0	100	0.00
33		Benzo(b)fluoranthene	1.000	1.063	-6.3	109	0.00
34		Benzo(k)fluoranthene	1.000	0.970	3.0	93	0.00
35	C	Benzo(a)pyrene	1.000	1.011	-1.1	101	-0.01
36		Dibenzo(a,h)anthracene	1.000	1.014	-1.4	100	-0.01
37		Benzo(g,h,i)perylene	1.000	1.011	-1.1	100	-0.01

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