Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM052015\
Data File : BM001365.D
Acq On : 21 May 2015 18:09
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 22 03:54:20 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 25 10:49:04 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	90	0.00
2		1,4-Dioxane	1.000	1.323	-32.3#	101	0.00
3		n-Nitrosodimethylamine	1.000	1.076	-7.6	91	0.00
4	S	2-Fluorophenol	1.000	1.028	-2.8	94	0.00
5	S	Phenol-d6	1.000	1.040	-4.0	95	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	90	0.00
7	S	Nitrobenzene-d5	1.000	1.031	-3.1	96	0.00
8		Nitrobenzene	1.000	0.959	4.1	96	0.00
9		Naphthalene	1.000	1.039	-3.9	93	0.00
10		Hexachlorobutadiene	1.000	1.046	-4.6	95	0.00
11		2-Methylnaphthalene	1.000	1.039	-3.9	94	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	91	0.00
13	S	2,4,6-Tribromophenol	1.000	1.041	-4.1	99	0.00
14	S	2-Fluorobiphenyl	1.000	1.045	-4.5	95	0.00
15		Acenaphthylene	1.000	1.040	-4.0	97	0.00
16		Acenaphthene	1.000	1.023	-2.3	95	0.00
17		Fluorene	1.000	0.981	1.9	91	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	86	0.00
19		4-Bromophenyl-phenylether	1.000	1.181	-18.1	106	0.00
20		Hexachlorobenzene	1.000	1.019	-1.9	88	0.00
21		Pentachlorophenol	1.000	1.002	-0.2	109	0.00
22		Phenanthrene	1.000	1.041	-4.1	89	0.00
23		Anthracene	1.000	1.174	-17.4	95	0.00
24		Fluoranthene	1.000	1.155	-15.5	99	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	96	0.00
26		Pyrene	1.000	1.012	-1.2	98	0.00
27	S	Terphenyl-d14	1.000	1.034	-3.4	101	0.00
28		Benzo(a)anthracene	1.000	1.033	-3.3	105	0.00
29		Chrysene	1.000	1.044	-4.4	103	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.918	8.2	102	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.074	-7.4	111	0.00
32	I	Perylene-d12	5.000	5.000	0.0	103	0.00
33		Benzo(b)fluoranthene	1.000	1.036	-3.6	110	0.00
34		Benzo(k)fluoranthene	1.000	1.000	0.0	104	0.00
35	C	Benzo(a)pyrene	1.000	1.022	-2.2	110	0.00
36		Dibenzo(a,h)anthracene	1.000	1.026	-2.6	111	0.00
37		Benzo(g,h,i)perylene	1.000	1.023	-2.3	111	0.00

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