Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050515\
Data File : BM001219.D
Acq On : 05 May 2015 21:52
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: May 06 13:12:17 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 06 12:46:35 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	98	0.00
2		1,4-Dioxane	1.000	0.881	11.9	85	0.00
3		n-Nitrosodimethylamine	1.000	0.922	7.8	100	0.00
4	S	2-Fluorophenol	1.000	0.993	0.7	95	0.00
5	S	Phenol-d6	1.000	0.968	3.2	92	-0.02
6	I	Naphthalene-d8	5.000	5.000	0.0	97	0.00
7	S	Nitrobenzene-d5	1.000	1.022	-2.2	96	0.00
8		Nitrobenzene	1.000	0.968	3.2	97	0.00
9		Naphthalene	1.000	1.027	-2.7	98	0.00
10		Hexachlorobutadiene	1.000	1.015	-1.5	99	0.00
11		2-Methylnaphthalene	1.000	1.024	-2.4	97	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13	S	2,4,6-Tribromophenol	1.000	0.912	8.8	94	0.00
14	S	2-Fluorobiphenyl	1.000	1.029	-2.9	96	-0.01
15		Acenaphthylene	1.000	1.020	-2.0	96	0.00
16		Acenaphthene	1.000	1.019	-1.9	96	0.00
17		Fluorene	1.000	0.986	1.4	88	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	91	0.00
19		Hexachlorobenzene	1.000	1.106	-10.6	101	0.00
20		Pentachlorophenol	1.000	1.036	-3.6	91	0.00
21		Phenanthrene	1.000	0.997	0.3	89	0.00
22		Anthracene	1.000	0.964	3.6	90	0.00
23		Fluoranthene	1.000	1.086	-8.6	95	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	97	0.00
25		Pyrene	1.000	0.990	1.0	96	0.00
26	S	Terphenyl-d14	1.000	0.987	1.3	95	0.00
27		Benzo(a)anthracene	1.000	1.019	-1.9	100	0.00
28		Chrysene	1.000	1.023	-2.3	98	0.00
29		Indeno(1,2,3-cd)pyrene	1.000	1.195	-19.5	114	0.00
30	I	Perylene-d12	5.000	5.000	0.0	106	0.00
31		Benzo(b)fluoranthene	1.000	0.961	3.9	106	0.00
32		Benzo(k)fluoranthene	1.000	1.051	-5.1	104	0.00
33	C	Benzo(a)pyrene	1.000	1.030	-3.0	108	0.00
34		Dibenzo(a,h)anthracene	1.000	1.080	-8.0	113	0.00
35		Benzo(g,h,i)perylene	1.000	1.094	-9.4	115	0.00

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