Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112715\
Data File : BM003311.D
Acq On : 27 Nov 2015 19:09
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Nov 27 22:49:17 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM112715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 27 16:36:40 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	103	0.00
2		1,4-Dioxane	1.000	1.092	-9.2	108	-0.02
3		n-Nitrosodimethylamine	1.000	1.000	0.0	103	0.00
4	S	2-Fluorophenol	1.000	0.984	1.6	104	0.00
5	S	Phenol-d6	1.000	0.968	3.2	103	0.00
6		bis(2-Chloroethyl)ether	1.000	0.972	2.8	103	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	101	0.00
8	S	Nitrobenzene-d5	1.000	0.960	4.0	102	0.00
9		Nitrobenzene	1.000	0.949	5.1	102	0.00
10		Naphthalene	1.000	0.997	0.3	102	0.00
11		Hexachlorobutadiene	1.000	1.056	-5.6	110	0.00
12		2-Methylnaphthalene	1.000	0.992	0.8	102	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
14	S	2,4,6-Tribromophenol	1.000	0.965	3.5	102	0.00
15	S	2-Fluorobiphenyl	1.000	0.932	6.8	102	0.00
16		Acenaphthylene	1.000	0.963	3.7	101	0.00
17		Acenaphthene	1.000	0.865	13.5	101	0.00
18		Fluorene	1.000	0.951	4.9	100	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	102	0.00
20		4-Bromophenyl-phenylether	1.000	0.838	16.2	101	0.00
21		Hexachlorobenzene	1.000	0.892	10.8	100	0.00
22		Pentachlorophenol	1.000	1.061	-6.1	130	0.00
23		Phenanthrene	1.000	0.834	16.6	96	0.03
24		Anthracene	1.000	0.972	2.8	100	0.00
25		Fluoranthene	1.000	0.882	11.8	101	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	92	0.00
27		Benzidine	1.000	1.324	-32.4#	153	0.00
28		Pyrene	1.000	0.927	7.3	102	0.00
29	S	Terphenyl-d14	1.000	0.960	4.0	97	0.00
30		Benzo(a)anthracene	1.000	1.014	-1.4	105	0.00
31		3,3'-Dichlorobenzidine	1.000	0.931	6.9	108	0.02
32		Chrysene	1.000	0.983	1.7	102	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.063	-6.3	122	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.996	0.4	106	0.01
35	I	Perylene-d12	5.000	5.000	0.0	103	0.01
36		Benzo(b)fluoranthene	1.000	0.965	3.5	97	0.01
37		Benzo(k)fluoranthene	1.000	1.035	-3.5	112	0.00
38	C	Benzo(a)pyrene	1.000	0.966	3.4	104	0.01
39		Dibenzo(a,h)anthracene	1.000	1.051	-5.1	109	0.01
40		Benzo(g,h,i)perylene	1.000	0.989	1.1	105	0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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