Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
Data File : BM004115.D
Acq On : 22 Jan 2016 17:39
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 22 18:17:55 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 22 14:46:41 2016
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	96	0.00
2		1,4-Dioxane	1.000	0.961	3.9	97	0.00
3		n-Nitrosodimethylamine	1.000	0.955	4.5	100	-0.02
4	S	2-Fluorophenol	1.000	0.919	8.1	98	-0.02
5	S	Phenol-d6	1.000	0.911	8.9	98	0.00
6		bis(2-Chloroethyl)ether	1.000	0.917	8.3	97	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
8	S	Nitrobenzene-d5	1.000	0.910	9.0	99	0.00
9		Nitrobenzene	1.000	0.911	8.9	100	0.00
10		Naphthalene	1.000	0.921	7.9	98	0.00
11		Hexachlorobutadiene	1.000	0.931	6.9	99	0.00
12		2-Methylnaphthalene	1.000	0.925	7.5	99	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
14	S	2,4,6-Tribromophenol	1.000	0.901	9.9	92	0.00
15	S	2-Fluorobiphenyl	1.000	0.979	2.1	98	0.00
16		Acenaphthylene	1.000	0.935	6.5	96	0.02
17		Acenaphthene	1.000	0.921	7.9	96	0.00
18		Fluorene	1.000	0.962	3.8	99	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
20		4-Bromophenyl-phenylether	1.000	0.806	19.4	86	0.02
21		Hexachlorobenzene	1.000	0.826	17.4	90	0.02
22		Pentachlorophenol	1.000	0.790	21.0	95	0.00
23		Phenanthrene	1.000	0.846	15.4	97	0.02
24		Anthracene	1.000	0.820	18.0	92	0.00
25		Fluoranthene	1.000	0.853	14.7	96	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
27		Benzidine	1.000	0.641	35.9#	58	0.00
28		Pyrene	1.000	0.862	13.8	95	0.00
29	S	Terphenyl-d14	1.000	0.897	10.3	100	0.00
30		Benzo(a)anthracene	1.000	0.814	18.6	94	0.00
31		3,3'-Dichlorobenzidine	1.000	0.884	11.6	96	-0.01
32		Chrysene	1.000	0.844	15.6	97	-0.01
33		Bis(2-ethylhexyl)phthalate	1.000	1.090	-9.0	127	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	0.858	14.2	99	0.00
35	I	Perylene-d12	5.000	5.000	0.0	97	0.00
36		Benzo(b)fluoranthene	1.000	0.904	9.6	94	0.00
37		Benzo(k)fluoranthene	1.000	0.934	6.6	102	0.00
38	C	Benzo(a)pyrene	1.000	0.921	7.9	100	0.00
39		Dibenzo(a,h)anthracene	1.000	0.917	8.3	99	0.00
40		Benzo(g,h,i)perylene	1.000	0.901	9.9	98	0.00

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

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