Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM020216\
Data File : BM004171.D
Acq On : 02 Feb 2016 19:34
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Feb 03 01:15:51 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	138	0.00
2		1,4-Dioxane	1.000	0.823	17.7	120	0.00
3		n-Nitrosodimethylamine	1.000	0.918	8.2	138	-0.02
4	S	2-Fluorophenol	1.000	0.942	5.8	143	-0.02
5	S	Phenol-d6	1.000	1.010	-1.0	155	0.00
6		bis(2-Chloroethyl)ether	1.000	0.934	6.6	142	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	146	0.00
8	S	Nitrobenzene-d5	1.000	1.018	-1.8	165	0.00
9		Nitrobenzene	1.000	0.995	0.5	167	0.00
10		Naphthalene	1.000	0.897	10.3	143	0.00
11		Hexachlorobutadiene	1.000	0.903	9.7	144	0.00
12		2-Methylnaphthalene	1.000	0.910	9.0	146	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
14	S	2,4,6-Tribromophenol	1.000	0.988	1.2	168	0.00
15	S	2-Fluorobiphenyl	1.000	0.873	12.7	141	0.00
16		Acenaphthylene	1.000	0.963	3.7	158	0.00
17		Acenaphthene	1.000	0.823	17.7	137	0.00
18		Fluorene	1.000	0.833	16.7	138	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	129	0.00
20		4-Bromophenyl-phenylether	1.000	1.186	-18.6	157	0.00
21		Hexachlorobenzene	1.000	1.137	-13.7	154	0.00
22		Pentachlorophenol	1.000	1.337	-33.7#	240	0.00
23		Phenanthrene	1.000	1.115	-11.5	153	0.00
24		Anthracene	1.000	1.046	-4.6	145	0.00
25		Fluoranthene	1.000	1.014	-1.4	141	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	142	0.00
27		Benzidine	1.000	1.309	-30.9#	202	0.00
28		Pyrene	1.000	0.924	7.6	144	0.00
29	S	Terphenyl-d14	1.000	0.928	7.2	147	0.00
30		Benzo(a)anthracene	1.000	0.864	13.6	140	0.00
31		3,3'-Dichlorobenzidine	1.000	1.104	-10.4	173	0.00
32		Chrysene	1.000	0.885	11.5	143	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.229	-22.9	206	-0.02
34		Indeno(1,2,3-cd)pyrene	1.000	0.787	21.3	127	0.00
35	I	Perylene-d12	5.000	5.000	0.0	131	0.00
36		Benzo(b)fluoranthene	1.000	0.903	9.7	127	-0.01
37		Benzo(k)fluoranthene	1.000	0.985	1.5	146	-0.01
38	C	Benzo(a)pyrene	1.000	0.958	4.2	141	0.00
39		Dibenzo(a,h)anthracene	1.000	0.879	12.1	128	-0.01
40		Benzo(g,h,i)perylene	1.000	0.835	16.5	123	-0.01

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

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