Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM122915\
Data File : BM003825.D
Acq On : 28 Dec 2015 13:04
Operator : UM/SJ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 29 01:30:04 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 10:08:49 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	109	0.00
2		1,4-Dioxane	1.000	0.926	7.4	112	0.00
3		n-Nitrosodimethylamine	1.000	0.857	14.3	108	0.00
4	S	2-Fluorophenol	1.000	0.915	8.5	117	0.02
5	S	Phenol-d6	1.000	0.922	7.8	121	0.00
6		bis(2-Chloroethyl)ether	1.000	0.881	11.9	111	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	111	-0.02
8	S	Nitrobenzene-d5	1.000	0.873	12.7	116	0.00
9		Nitrobenzene	1.000	0.873	12.7	115	0.00
10		Naphthalene	1.000	0.915	8.5	115	0.00
11		Hexachlorobutadiene	1.000	0.930	7.0	117	0.00
12		2-Methylnaphthalene	1.000	0.918	8.2	117	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	114	0.00
14	S	2,4,6-Tribromophenol	1.000	0.866	13.4	119	0.00
15	S	2-Fluorobiphenyl	1.000	0.904	9.6	118	0.00
16		Acenaphthylene	1.000	0.902	9.8	125	0.00
17		Acenaphthene	1.000	0.910	9.0	112	0.00
18		Fluorene	1.000	0.885	11.5	117	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	117	0.00
20		4-Bromophenyl-phenylether	1.000	1.221	-22.1	112	0.00
21		Hexachlorobenzene	1.000	1.201	-20.1	112	0.00
22		Pentachlorophenol	1.000	0.649	35.1#	70	0.00
23		Phenanthrene	1.000	1.120	-12.0	113	0.00
24		Anthracene	1.000	0.914	8.6	115	0.00
25		Fluoranthene	1.000	0.894	10.6	109	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
27		Benzidine	1.000	1.369	-36.9#	142	0.00
28		Pyrene	1.000	1.060	-6.0	107	0.00
29	S	Terphenyl-d14	1.000	1.017	-1.7	101	0.00
30		Benzo(a)anthracene	1.000	0.914	8.6	102	0.00
31		3,3'-Dichlorobenzidine	1.000	1.090	-9.0	109	0.00
32		Chrysene	1.000	0.871	12.9	85	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.144	-14.4	126	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.171	-17.1	117	0.01
35	I	Perylene-d12	5.000	5.000	0.0	95	0.01
36		Benzo(b)fluoranthene	1.000	0.860	14.0	90	0.00
37		Benzo(k)fluoranthene	1.000	0.918	8.2	104	0.00
38	C	Benzo(a)pyrene	1.000	0.943	5.7	105	0.00
39		Dibenzo(a,h)anthracene	1.000	1.086	-8.6	117	0.00
40		Benzo(g,h,i)perylene	1.000	1.115	-11.5	119	0.00

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

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