Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120215\
Data File : BM003366.D
Acq On : 02 Dec 2015 21:11
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 03 00:44:55 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 16:19:00 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	66	0.00
2		1,4-Dioxane	1.000	0.937	6.3	69	0.00
3		n-Nitrosodimethylamine	1.000	0.947	5.3	65	0.00
4	S	2-Fluorophenol	1.000	0.867	13.3	64	0.00
5	S	Phenol-d6	1.000	0.885	11.5	62	0.00
6		bis(2-Chloroethyl)ether	1.000	0.958	4.2	66	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	64	0.00
8	S	Nitrobenzene-d5	1.000	0.894	10.6	62	0.00
9		Nitrobenzene	1.000	0.896	10.4	61	0.00
10		Naphthalene	1.000	0.949	5.1	66	0.00
11		Hexachlorobutadiene	1.000	0.979	2.1	68	0.00
12		2-Methylnaphthalene	1.000	0.944	5.6	63	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	58	0.00
14	S	2,4,6-Tribromophenol	1.000	1.060	-6.0	62	0.00
15	S	2-Fluorobiphenyl	1.000	0.998	0.2	64	0.00
16		Acenaphthylene	1.000	0.919	8.1	54	0.00
17		Acenaphthene	1.000	1.009	-0.9	65	-0.02
18		Fluorene	1.000	1.122	-12.2	69	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	55	0.00
20		4-Bromophenyl-phenylether	1.000	1.175	-17.5	75	-0.02
21		Hexachlorobenzene	1.000	1.323	-32.3#	80	0.00
22		Pentachlorophenol	1.000	0.864	13.6	47	-0.02
23		Phenanthrene	1.000	1.155	-15.5	72	-0.02
24		Anthracene	1.000	0.993	0.7	56	0.00
25		Fluoranthene	1.000	1.083	-8.3	59	-0.01
26	I	Chrysene-d12	5.000	5.000	0.0	67	-0.01
27		Benzidine	1.000	0.772	22.8	49	0.00
28		Pyrene	1.000	0.770	23.0	57	0.00
29	S	Terphenyl-d14	1.000	0.832	16.8	60	0.00
30		Benzo(a)anthracene	1.000	0.857	14.3	60	-0.01
31		3,3'-Dichlorobenzidine	1.000	0.608	39.2#	43	0.00
32		Chrysene	1.000	0.731	26.9#	52	0.01
33		Bis(2-ethylhexyl)phthalate	1.000	0.604	39.6#	32	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	0.636	36.4#	55	0.00
35	I	Perylene-d12	5.000	5.000	0.0	54	0.00
36		Benzo(b)fluoranthene	1.000	0.965	3.5	54	-0.01
37		Benzo(k)fluoranthene	1.000	0.911	8.9	51	0.00
38	C	Benzo(a)pyrene	1.000	0.846	15.4	49	0.00
39		Dibenzo(a,h)anthracene	1.000	0.847	15.3	57	0.00
40		Benzo(g,h,i)perylene	1.000	0.843	15.7	58	0.00

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

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