Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120415\
Data File : BM003377.D
Acq On : 03 Dec 2015 19:03
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 04 00:19:33 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	63	0.00
2		1,4-Dioxane	1.000	0.914	8.6	65	0.00
3		n-Nitrosodimethylamine	1.000	0.982	1.8	64	0.00
4	S	2-Fluorophenol	1.000	0.939	6.1	67	0.00
5	S	Phenol-d6	1.000	0.964	3.6	65	0.00
6		bis(2-Chloroethyl)ether	1.000	0.960	4.0	63	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	62	0.00
8	S	Nitrobenzene-d5	1.000	0.967	3.3	65	0.00
9		Nitrobenzene	1.000	0.965	3.5	64	0.00
10		Naphthalene	1.000	0.951	4.9	64	0.00
11		Hexachlorobutadiene	1.000	0.988	1.2	67	0.00
12		2-Methylnaphthalene	1.000	0.960	4.0	63	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	61	0.00
14	S	2,4,6-Tribromophenol	1.000	1.065	-6.5	65	0.00
15	S	2-Fluorobiphenyl	1.000	0.929	7.1	63	0.00
16		Acenaphthylene	1.000	0.973	2.7	61	0.00
17		Acenaphthene	1.000	0.908	9.2	62	0.00
18		Fluorene	1.000	0.981	1.9	63	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	62	0.00
20		4-Bromophenyl-phenylether	1.000	0.901	9.9	64	-0.03
21		Hexachlorobenzene	1.000	0.980	2.0	66	0.00
22		Pentachlorophenol	1.000	0.990	1.0	64	0.00
23		Phenanthrene	1.000	1.031	-3.1	72	0.08
24		Anthracene	1.000	1.000	0.0	63	0.00
25		Fluoranthene	1.000	1.027	-2.7	63	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	64	0.00
27		Benzidine	1.000	0.595	40.5#	26	0.00
28		Pyrene	1.000	0.884	11.6	62	0.00
29	S	Terphenyl-d14	1.000	0.949	5.1	65	0.00
30		Benzo(a)anthracene	1.000	0.954	4.6	64	0.00
31		3,3'-Dichlorobenzidine	1.000	0.890	11.0	62	0.00
32		Chrysene	1.000	0.988	1.2	67	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.039	-3.9	68	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.797	20.3	66	0.00
35	I	Perylene-d12	5.000	5.000	0.0	66	0.00
36		Benzo(b)fluoranthene	1.000	0.984	1.6	68	0.00
37		Benzo(k)fluoranthene	1.000	0.948	5.2	66	0.00
38	C	Benzo(a)pyrene	1.000	0.921	7.9	66	0.00
39		Dibenzo(a,h)anthracene	1.000	0.799	20.1	66	0.00
40		Benzo(g,h,i)perylene	1.000	0.795	20.5	68	0.00

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

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