Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101615\
Data File : BM002937.D
Acq On : 16 Oct 2015 16:42
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 16 23:25:32 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM101515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 16 22:59:32 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	110	0.00
2		1,4-Dioxane	1.000	0.941	5.9	106	0.00
3		n-Nitrosodimethylamine	1.000	1.073	-7.3	115	-0.03
4	S	2-Fluorophenol	1.000	1.046	-4.6	118	-0.03
5	S	Phenol-d6	1.000	1.085	-8.5	118	-0.02
6		bis(2-Chloroethyl)ether	1.000	1.030	-3.0	111	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	112	-0.01
8	S	Nitrobenzene-d5	1.000	1.053	-5.3	118	-0.03
9		Nitrobenzene	1.000	1.058	-5.8	118	-0.03
10		Naphthalene	1.000	0.963	3.7	110	-0.01
11		Hexachlorobutadiene	1.000	0.975	2.5	111	-0.01
12		2-Methylnaphthalene	1.000	0.983	1.7	112	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	109	0.00
14	S	2,4,6-Tribromophenol	1.000	1.030	-3.0	123	0.00
15	S	2-Fluorobiphenyl	1.000	0.999	0.1	111	-0.01
16		Acenaphthylene	1.000	1.056	-5.6	117	0.00
17		Acenaphthene	1.000	0.978	2.2	111	0.00
18		Fluorene	1.000	1.005	-0.5	111	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	116	0.00
20		4-Bromophenyl-phenylether	1.000	1.021	-2.1	122	0.00
21		Hexachlorobenzene	1.000	1.048	-4.8	98	0.00
22		Pentachlorophenol	1.000	1.689	-68.9#	225	-0.03
23		Phenanthrene	1.000	0.884	11.6	104	0.02
24		Anthracene	1.000	0.933	6.7	108	0.00
25		Fluoranthene	1.000	0.938	6.2	112	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
27		Benzidine	1.000	0.815	18.5	87	0.00
28		Pyrene	1.000	1.137	-13.7	109	0.00
29	S	Terphenyl-d14	1.000	1.109	-10.9	106	0.00
30		Benzo(a)anthracene	1.000	1.043	-4.3	96	0.00
31		3,3'-Dichlorobenzidine	1.000	1.081	-8.1	108	0.00
32		Chrysene	1.000	1.003	-0.3	104	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.339	-33.9#	134	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.800	20.0	82	0.00
35	I	Perylene-d12	5.000	5.000	0.0	90	0.00
36		Benzo(b)fluoranthene	1.000	1.052	-5.2	95	0.00
37		Benzo(k)fluoranthene	1.000	1.018	-1.8	95	-0.01
38	C	Benzo(a)pyrene	1.000	1.027	-2.7	95	0.00
39		Dibenzo(a,h)anthracene	1.000	0.881	11.9	83	-0.01
40		Benzo(g,h,i)perylene	1.000	0.851	14.9	80	-0.01

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

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