Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101515\
Data File : BM002928.D
Acq On : 15 Oct 2015 23:34
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 16 07:32:31 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 06:59:27 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	101	0.00
2		1,4-Dioxane	1.000	0.943	5.7	98	0.00
3		n-Nitrosodimethylamine	1.000	1.050	-5.0	103	-0.03
4	S	2-Fluorophenol	1.000	1.019	-1.9	105	-0.03
5	S	Phenol-d6	1.000	1.068	-6.8	106	-0.02
6		bis(2-Chloroethyl)ether	1.000	1.021	-2.1	101	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	102	-0.01
8	S	Nitrobenzene-d5	1.000	1.033	-3.3	105	-0.03
9		Nitrobenzene	1.000	1.036	-3.6	105	-0.03
10		Naphthalene	1.000	0.966	3.4	100	-0.01
11		Hexachlorobutadiene	1.000	0.982	1.8	102	-0.01
12		2-Methylnaphthalene	1.000	0.973	2.7	101	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	99	0.00
14	S	2,4,6-Tribromophenol	1.000	1.056	-5.6	116	0.00
15	S	2-Fluorobiphenyl	1.000	0.986	1.4	100	-0.01
16		Acenaphthylene	1.000	1.029	-2.9	104	0.00
17		Acenaphthene	1.000	0.954	4.6	99	0.00
18		Fluorene	1.000	0.997	0.3	101	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	106	0.00
20		4-Bromophenyl-phenylether	1.000	1.017	-1.7	111	0.00
21		Hexachlorobenzene	1.000	1.052	-5.2	90	0.00
22		Pentachlorophenol	1.000	1.470	-47.0#	175	-0.03
23		Phenanthrene	1.000	0.875	12.5	94	0.02
24		Anthracene	1.000	0.906	9.4	96	0.00
25		Fluoranthene	1.000	0.913	8.7	100	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	95	0.00
27		Benzidine	1.000	1.009	-0.9	113	0.00
28		Pyrene	1.000	1.021	-2.1	98	0.00
29	S	Terphenyl-d14	1.000	1.031	-3.1	99	0.00
30		Benzo(a)anthracene	1.000	1.080	-8.0	100	0.00
31		3,3'-Dichlorobenzidine	1.000	1.114	-11.4	112	0.00
32		Chrysene	1.000	0.899	10.1	93	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.248	-24.8	123	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.869	13.1	90	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	93	0.00
36		Benzo(b)fluoranthene	1.000	1.060	-6.0	98	-0.01
37		Benzo(k)fluoranthene	1.000	0.960	4.0	92	-0.01
38	C	Benzo(a)pyrene	1.000	1.005	-0.5	96	-0.01
39		Dibenzo(a,h)anthracene	1.000	0.934	6.6	91	-0.01
40		Benzo(g,h,i)perylene	1.000	0.898	10.2	88	-0.01

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

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