Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM100915\
Data File : BM002889.D
Acq On : 09 Oct 2015 17:47
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 12 05:15:35 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM100815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 09 06:47:52 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	152	-0.02
2		1,4-Dioxane	1.000	0.956	4.4	141	0.00
3		n-Nitrosodimethylamine	1.000	0.949	5.1	140	0.00
4	S	2-Fluorophenol	1.000	0.870	13.0	134	0.00
5	S	Phenol-d6	1.000	0.910	9.0	143	0.00
6		bis(2-Chloroethyl)ether	1.000	0.955	4.5	150	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	150	-0.01
8	S	Nitrobenzene-d5	1.000	0.953	4.7	145	0.00
9		Nitrobenzene	1.000	0.953	4.7	146	0.00
10		Naphthalene	1.000	0.961	3.9	148	-0.01
11		Hexachlorobutadiene	1.000	0.976	2.4	151	0.00
12		2-Methylnaphthalene	1.000	0.959	4.1	147	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	149	0.00
14	S	2,4,6-Tribromophenol	1.000	0.827	17.3	133	0.00
15	S	2-Fluorobiphenyl	1.000	0.986	1.4	147	0.00
16		Acenaphthylene	1.000	0.964	3.6	145	0.00
17		Acenaphthene	1.000	0.938	6.2	146	0.00
18		Fluorene	1.000	0.949	5.1	145	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	131	0.02
20		4-Bromophenyl-phenylether	1.000	0.913	8.7	117	0.00
21		Hexachlorobenzene	1.000	1.117	-11.7	149	0.00
22		Pentachlorophenol	1.000	0.771	22.9	89	0.02
23		Phenanthrene	1.000	0.998	0.2	134	0.00
24		Anthracene	1.000	1.086	-8.6	149	0.00
25		Fluoranthene	1.000	0.925	7.5	130	0.01
26	I	Chrysene-d12	5.000	5.000	0.0	126	0.00
27		Benzidine	1.000	0.763	23.7	101	0.00
28		Pyrene	1.000	1.032	-3.2	128	0.01
29	S	Terphenyl-d14	1.000	1.056	-5.6	136	0.00
30		Benzo(a)anthracene	1.000	0.901	9.9	122	0.00
31		3,3'-Dichlorobenzidine	1.000	0.814	18.6	107	0.00
32		Chrysene	1.000	0.965	3.5	123	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.873	12.7	93	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.681	31.9#	88	0.00
35	I	Perylene-d12	5.000	5.000	0.0	101	0.00
36		Benzo(b)fluoranthene	1.000	0.987	1.3	98	0.00
37		Benzo(k)fluoranthene	1.000	1.030	-3.0	111	0.00
38	C	Benzo(a)pyrene	1.000	0.944	5.6	98	0.00
39		Dibenzo(a,h)anthracene	1.000	0.849	15.1	88	0.00
40		Benzo(g,h,i)perylene	1.000	0.848	15.2	88	0.00

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

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