Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004490.D
Acq On : 01 Mar 2016 03:47
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Mar 01 07:29:49 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM022916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 01 00:52:28 2016
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	84	0.00
2		1,4-Dioxane	1.000	0.939	6.1	80	0.00
3		n-Nitrosodimethylamine	1.000	0.871	12.9	78	0.00
4	S	2-Fluorophenol	1.000	0.877	12.3	79	0.00
5	S	Phenol-d6	1.000	0.942	5.8	83	0.02
6		bis(2-Chloroethyl)ether	1.000	0.936	6.4	81	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	88	0.00
8	S	Nitrobenzene-d5	1.000	0.911	8.9	85	0.00
9		Nitrobenzene	1.000	0.904	9.6	83	0.00
10		Naphthalene	1.000	0.938	6.2	86	0.00
11		Hexachlorobutadiene	1.000	0.966	3.4	87	-0.01
12		2-Methylnaphthalene	1.000	0.995	0.5	92	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
14	S	2,4,6-Tribromophenol	1.000	0.862	13.8	89	0.00
15	S	2-Fluorobiphenyl	1.000	0.963	3.7	93	0.00
16		Acenaphthylene	1.000	0.944	5.6	91	0.00
17		Acenaphthene	1.000	0.972	2.8	100	0.00
18		Fluorene	1.000	1.005	-0.5	98	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	87	0.00
20		4-Bromophenyl-phenylether	1.000	1.078	-7.8	109	0.00
21		Hexachlorobenzene	1.000	1.088	-8.8	114	0.00
22		Pentachlorophenol	1.000	1.101	-10.1	122	0.00
23		Phenanthrene	1.000	1.031	-3.1	106	0.00
24		Anthracene	1.000	0.887	11.3	88	0.02
25		Fluoranthene	1.000	0.957	4.3	98	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	79	-0.01
27		Benzidine	1.000	0.797	20.3	72	0.00
28		Pyrene	1.000	1.078	-7.8	98	0.00
29	S	Terphenyl-d14	1.000	1.041	-4.1	95	0.00
30		Benzo(a)anthracene	1.000	0.965	3.5	86	-0.01
31		3,3'-Dichlorobenzidine	1.000	1.034	-3.4	99	0.00
32		Chrysene	1.000	0.834	16.6	75	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.909	9.1	69	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	1.042	-4.2	92	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	93	-0.01
36		Benzo(b)fluoranthene	1.000	1.137	-13.7	107	-0.01
37		Benzo(k)fluoranthene	1.000	1.010	-1.0	99	0.00
38	C	Benzo(a)pyrene	1.000	0.989	1.1	95	0.00
39		Dibenzo(a,h)anthracene	1.000	0.956	4.4	93	-0.01
40		Benzo(g,h,i)perylene	1.000	0.960	4.0	93	-0.01

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

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