Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\DATA\BM022916\
Data File : BM004500.D
Acq On : 01 Mar 2016 10:29
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Mar 01 11:57:37 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	90	0.00
2		1,4-Dioxane	1.000	0.996	0.4	90	0.00
3		n-Nitrosodimethylamine	1.000	0.910	9.0	88	0.00
4	S	2-Fluorophenol	1.000	0.917	8.3	89	0.02
5	S	Phenol-d6	1.000	0.986	1.4	94	0.02
6		bis(2-Chloroethyl)ether	1.000	0.987	1.3	92	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
8	S	Nitrobenzene-d5	1.000	0.938	6.2	93	0.00
9		Nitrobenzene	1.000	0.927	7.3	91	0.00
10		Naphthalene	1.000	0.971	2.9	95	0.00
11		Hexachlorobutadiene	1.000	1.000	0.0	96	-0.01
12		2-Methylnaphthalene	1.000	1.027	-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	0.890	11.0	99	0.00
15	S	2-Fluorobiphenyl	1.000	0.993	0.7	104	0.00
16		Acenaphthylene	1.000	0.949	5.1	99	0.00
17		Acenaphthene	1.000	1.012	-1.2	112	0.00
18		Fluorene	1.000	1.030	-3.0	108	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
20		4-Bromophenyl-phenylether	1.000	1.184	-18.4	123	0.00
21		Hexachlorobenzene	1.000	1.186	-18.6	128	0.00
22		Pentachlorophenol	1.000	1.153	-15.3	132	0.00
23		Phenanthrene	1.000	1.134	-13.4	120	0.00
24		Anthracene	1.000	0.981	1.9	100	0.02
25		Fluoranthene	1.000	1.039	-3.9	109	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	84	0.01
27		Benzidine	1.000	1.315	-31.5#	161	0.00
28		Pyrene	1.000	1.118	-11.8	108	0.00
29	S	Terphenyl-d14	1.000	1.110	-11.0	107	0.00
30		Benzo(a)anthracene	1.000	0.960	4.0	91	0.01
31		3,3'-Dichlorobenzidine	1.000	1.510	-51.0#	161	0.00
32		Chrysene	1.000	0.949	5.1	89	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.831	16.9	67	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	1.087	-8.7	102	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	100	0.00
36		Benzo(b)fluoranthene	1.000	1.043	-4.3	104	0.00
37		Benzo(k)fluoranthene	1.000	0.989	1.1	104	0.00
38	C	Benzo(a)pyrene	1.000	1.011	-1.1	103	0.00
39		Dibenzo(a,h)anthracene	1.000	0.985	1.5	102	0.00
40		Benzo(g,h,i)perylene	1.000	0.987	1.3	102	-0.01

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

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