Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
Data File : BM004107.D
Acq On : 21 Jan 2016 19:52
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDICC002

Quant Time: Jan 22 01:22:07 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM010216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 13:33:21 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	54	0.00
2		1,4-Dioxane	1.000	1.009	-0.9	54	0.00
3		n-Nitrosodimethylamine	1.000	0.884	11.6	48	-0.02
4	S	2-Fluorophenol	1.000	0.951	4.9	55	-0.02
5	S	Phenol-d6	1.000	0.934	6.6	52	0.00
6		bis(2-Chloroethyl)ether	1.000	0.952	4.8	50	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	54	-0.02
8	S	Nitrobenzene-d5	1.000	0.921	7.9	50	-0.01
9		Nitrobenzene	1.000	0.911	8.9	50	-0.01
10		Naphthalene	1.000	0.985	1.5	54	-0.02
11		Hexachlorobutadiene	1.000	0.991	0.9	55	0.00
12		2-Methylnaphthalene	1.000	1.014	-1.4	55	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	60	-0.02
14	S	2,4,6-Tribromophenol	1.000	1.097	-9.7	73	0.00
15	S	2-Fluorobiphenyl	1.000	0.944	5.6	57	-0.01
16		Acenaphthylene	1.000	1.002	-0.2	59	-0.02
17		Acenaphthene	1.000	1.017	-1.7	65	0.00
18		Fluorene	1.000	1.046	-4.6	62	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	63	0.00
20		4-Bromophenyl-phenylether	1.000	1.017	-1.7	68	-0.03
21		Hexachlorobenzene	1.000	1.043	-4.3	73	-0.03
22		Pentachlorophenol	1.000	1.185	-18.5	84	0.00
23		Phenanthrene	1.000	1.052	-5.2	65	-0.02
24		Anthracene	1.000	1.109	-10.9	76	0.00
25		Fluoranthene	1.000	1.137	-13.7	75	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	103	0.00
27		Benzidine	1.000	0.595	40.5#	52	0.00
28		Pyrene	1.000	0.786	21.4	81	0.00
29	S	Terphenyl-d14	1.000	0.775	22.5	81	-0.02
30		Benzo(a)anthracene	1.000	1.021	-2.1	108	0.00
31		3,3'-Dichlorobenzidine	1.000	0.962	3.8	105	0.00
32		Chrysene	1.000	0.888	11.2	90	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.929	7.1	108	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.407	-40.7#	145	-0.01
35	I	Perylene-d12	5.000	5.000	0.0	132	-0.01
36		Benzo(b)fluoranthene	1.000	0.991	0.9	140	0.00
37		Benzo(k)fluoranthene	1.000	0.870	13.0	117	0.00
38	C	Benzo(a)pyrene	1.000	0.995	0.5	139	-0.01
39		Dibenzo(a,h)anthracene	1.000	1.086	-8.6	144	-0.01
40		Benzo(g,h,i)perylene	1.000	1.072	-7.2	143	0.00

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

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