Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
Data File : BM004102.D
Acq On : 21 Jan 2016 16:53
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDICC0.1

Quant Time: Jan 22 01:26:24 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	53	0.00
2		1,4-Dioxane	1.000	1.012	-1.2	54	0.00
3		n-Nitrosodimethylamine	1.000	0.885	11.5	48	-0.02
4	S	2-Fluorophenol	1.000	0.960	4.0	55	-0.02
5	S	Phenol-d6	1.000	0.940	6.0	52	0.00
6		bis(2-Chloroethyl)ether	1.000	1.120	-12.0	59	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	54	-0.02
8	S	Nitrobenzene-d5	1.000	0.927	7.3	51	-0.01
9		Nitrobenzene	1.000	0.918	8.2	51	-0.01
10		Naphthalene	1.000	0.989	1.1	54	-0.02
11		Hexachlorobutadiene	1.000	0.990	1.0	55	0.00
12		2-Methylnaphthalene	1.000	1.018	-1.8	55	-0.01
13	I	Acenaphthene-d10	5.000	5.000	0.0	65	-0.02
14	S	2,4,6-Tribromophenol	1.000	1.016	-1.6	72	0.00
15	S	2-Fluorobiphenyl	1.000	0.863	13.7	57	-0.01
16		Acenaphthylene	1.000	0.974	2.6	63	-0.02
17		Acenaphthene	1.000	0.847	15.3	59	0.00
18		Fluorene	1.000	0.865	13.5	56	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	50	0.00
20		4-Bromophenyl-phenylether	1.000	1.667	-66.7#	89	-0.02
21		Hexachlorobenzene	1.000	1.580	-58.0#	88	-0.02
22		Pentachlorophenol	1.000	1.405	-40.5#	82	0.00
23		Phenanthrene	1.000	1.721	-72.1#	82	-0.02
24		Anthracene	1.000	1.270	-27.0#	69	0.00
25		Fluoranthene	1.000	1.466	-46.6#	77	-0.01
26	I	Chrysene-d12	5.000	5.000	0.0	109	0.01
27		Benzidine	1.000	0.650	35.0#	62	0.00
28		Pyrene	1.000	0.734	26.6#	79	0.00
29	S	Terphenyl-d14	1.000	0.777	22.3	85	-0.01
30		Benzo(a)anthracene	1.000	0.979	2.1	110	0.01
31		3,3'-Dichlorobenzidine	1.000	1.079	-7.9	128	0.00
32		Chrysene	1.000	0.901	9.9	97	-0.01
33		Bis(2-ethylhexyl)phthalate	1.000	1.126	-12.6	153	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	1.374	-37.4#	149	0.00
35	I	Perylene-d12	5.000	5.000	0.0	137	-0.01
36		Benzo(b)fluoranthene	1.000	0.966	3.4	141	0.00
37		Benzo(k)fluoranthene	1.000	0.884	11.6	124	0.00
38	C	Benzo(a)pyrene	1.000	0.994	0.6	144	0.00
39		Dibenzo(a,h)anthracene	1.000	1.074	-7.4	147	-0.01
40		Benzo(g,h,i)perylene	1.000	1.066	-6.6	147	0.00

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

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