Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010816\
Data File : BM003929.D
Acq On : 08 Jan 2016 19:09
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 09 00:54:45 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	82	0.00
2		1,4-Dioxane	1.000	0.990	1.0	82	0.00
3		n-Nitrosodimethylamine	1.000	0.872	12.8	72	0.00
4	S	2-Fluorophenol	1.000	0.907	9.3	80	0.00
5	S	Phenol-d6	1.000	0.924	7.6	78	0.00
6		bis(2-Chloroethyl)ether	1.000	0.967	3.3	78	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	82	0.00
8	S	Nitrobenzene-d5	1.000	0.854	14.6	71	0.00
9		Nitrobenzene	1.000	0.909	9.1	76	0.00
10		Naphthalene	1.000	0.964	3.6	79	0.00
11		Hexachlorobutadiene	1.000	0.960	4.0	80	0.01
12		2-Methylnaphthalene	1.000	0.913	8.7	75	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	82	-0.02
14	S	2,4,6-Tribromophenol	1.000	0.918	8.2	80	0.00
15	S	2-Fluorobiphenyl	1.000	0.902	9.8	75	0.00
16		Acenaphthylene	1.000	0.925	7.5	76	-0.02
17		Acenaphthene	1.000	0.989	1.1	87	0.00
18		Fluorene	1.000	0.962	3.8	78	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	75	0.00
20		4-Bromophenyl-phenylether	1.000	1.127	-12.7	90	-0.02
21		Hexachlorobenzene	1.000	1.178	-17.8	98	-0.02
22		Pentachlorophenol	1.000	0.942	5.8	74	0.00
23		Phenanthrene	1.000	1.182	-18.2	86	-0.02
24		Anthracene	1.000	1.089	-8.9	88	0.00
25		Fluoranthene	1.000	1.029	-2.9	80	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	97	0.01
27		Benzidine	1.000	0.124	87.6#	0	-19.33#
28		Pyrene	1.000	0.865	13.5	83	0.00
29	S	Terphenyl-d14	1.000	0.883	11.7	86	-0.01
30		Benzo(a)anthracene	1.000	0.877	12.3	87	0.01
31		3,3'-Dichlorobenzidine	1.000	0.675	32.5#	59	0.00
32		Chrysene	1.000	0.932	6.8	89	-0.01
33		Bis(2-ethylhexyl)phthalate	1.000	0.965	3.5	108	-0.01
34		Indeno(1,2,3-cd)pyrene	1.000	0.854	14.6	83	0.00
35	I	Perylene-d12	5.000	5.000	0.0	95	-0.01
36		Benzo(b)fluoranthene	1.000	0.970	3.0	99	0.00
37		Benzo(k)fluoranthene	1.000	0.892	10.8	87	0.00
38	C	Benzo(a)pyrene	1.000	0.921	7.9	93	0.00
39		Dibenzo(a,h)anthracene	1.000	0.864	13.6	82	-0.01
40		Benzo(g,h,i)perylene	1.000	0.873	12.7	84	0.00

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

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