Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM022816\
Data File : BM004473.D
Acq On : 29 Feb 2016 09:38
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Feb 29 10:41:58 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM022816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 29 05:57:18 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	95	0.00
2		1,4-Dioxane	1.000	0.955	4.5	102	0.00
3		n-Nitrosodimethylamine	1.000	0.855	14.5	94	0.00
4	S	2-Fluorophenol	1.000	0.793	20.7	86	0.02
5	S	Phenol-d6	1.000	0.785	21.5	93	0.00
6		bis(2-Chloroethyl)ether	1.000	0.809	19.1	87	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
8	S	Nitrobenzene-d5	1.000	0.742	25.8#	88	0.00
9		Nitrobenzene	1.000	0.742	25.8#	86	0.00
10		Hexachlorobutadiene	1.000	0.808	19.2	92	0.00
11		2-Methylnaphthalene	1.000	0.876	12.4	101	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
13	S	2,4,6-Tribromophenol	1.000	0.596	40.4#	73	0.00
14	S	2-Fluorobiphenyl	1.000	0.755	24.5	100	0.00
15	I	Phenanthrene-d10	5.000	5.000	0.0	98	0.00
16		Hexachlorobenzene	1.000	0.926	7.4	98	0.00
17		Pentachlorophenol	1.000	0.767	23.3	90	0.00
18		Fluoranthene	1.000	0.891	10.9	104	0.00
19	I	Chrysene-d12	5.000	5.000	0.0	105	0.00
20		Benzidine	1.000	1.025	-2.5	147	0.00
21		Pyrene	1.000	0.877	12.3	105	0.00
22	S	Terphenyl-d14	1.000	0.871	12.9	105	0.00
23		Benzo(a)anthracene	1.000	0.790	21.0	106	0.00
24		3,3'-Dichlorobenzidine	1.000	0.902	9.8	99	0.00
25		Chrysene	1.000	0.588	41.2#	90	0.02
26		Bis(2-ethylhexyl)phthalate	1.000	0.724	27.6#	89	0.00
27		Indeno(1,2,3-cd)pyrene	1.000	0.865	13.5	106	0.00
28	I	Perylene-d12	5.000	5.000	0.0	108	0.00
29		Benzo(b)fluoranthene	1.000	0.872	12.8	104	0.00
30		Benzo(k)fluoranthene	1.000	0.825	17.5	110	0.00
31	C	Benzo(a)pyrene	1.000	0.879	12.1	107	0.00
32		Dibenzo(a,h)anthracene	1.000	0.829	17.1	105	0.00
33		Benzo(g,h,i)perylene	1.000	0.844	15.6	107	0.00

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