Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000830.D
Acq On : 03 Apr 2015 06:31
Operator : TP/IZ
Sample : SSTDCCC001
Misc : MDL-WATER- BLK
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 03 18:40:23 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 03 18:16:53 2015
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	1.179	-17.9	107	0.00
3		n-Nitrosodimethylamine	1.000	1.054	-5.4	100	0.00
4	S	2-Fluorophenol	1.000	1.097	-9.7	104	0.00
5	S	Phenol-d6	1.000	1.066	-6.6	109	0.00
6		2-Chlorophenol	1.000	1.099	-9.9	111	0.00
7	C	Phenol	1.000	1.030	-3.0	100	0.00
8		bis(2-Chloroethyl)ether	1.000	1.050	-5.0	109	0.00
9		2-Methylphenol	1.000	1.037	-3.7	103	0.00
10		3+4-Methylphenols	1.000	1.025	-2.5	103	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	101	0.00
12	S	Nitrobenzene-d5	1.000	0.985	1.5	103	0.00
13		Nitrobenzene	1.000	0.962	3.8	100	0.00
14		2-Nitrophenol	1.000	1.021	-2.1	103	0.00
15		2,4-Dimethylphenol	1.000	1.039	-3.9	102	0.00
16	C	2,4-Dichlorophenol	1.000	1.002	-0.2	103	0.00
17	C	Hexachlorobutadiene	1.000	1.042	-4.2	103	-0.03
18		4-Chloro-3-methylphenol	1.000	1.031	-3.1	105	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
20	S	2,4,6-Tribromophenol	1.000	0.898	10.2	99	0.00
21		2,4,6-Trichlorophenol	1.000	0.897	10.3	94	0.00
22		2,4,5-Trichlorophenol	1.000	1.043	-4.3	102	0.00
23	S	2-Fluorobiphenyl	1.000	1.011	-1.1	103	0.00
24	P	2,4-Dinitrophenol	1.000	0.285	71.5#	65	0.00
25		4-Nitrophenol	1.000	0.917	8.3	95	0.00
26	I	Phenanthrene-d10	5.000	5.000	0.0	106	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.809	19.1	82	0.03
28		Hexachlorobenzene	1.000	0.944	5.6	100	0.00
29	C	Pentachlorophenol	1.000	0.817	18.3	75	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	103	0.00
31		Benzidine	1.000	1.674	-67.4#	1187	-0.02
32	S	Terphenyl-d14	1.000	1.019	-1.9	103	0.00
33		3,3'-Dichlorobenzidine	1.000	0.990	1.0	138	0.00
34	I	Perylene-d12	5.000	5.000	0.0	104	0.00

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