Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040315\
Data File : BM000841.D
Acq On : 03 Apr 2015 18:48
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 04 07:18:45 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	1.169	-16.9	105	0.00
3		n-Nitrosodimethylamine	1.000	1.119	-11.9	106	-0.02
4	S	2-Fluorophenol	1.000	1.171	-17.1	110	0.00
5	S	Phenol-d6	1.000	1.149	-14.9	119	0.00
6		2-Chlorophenol	1.000	1.225	-22.5	126	0.00
7	C	Phenol	1.000	1.063	-6.3	103	0.00
8		bis(2-Chloroethyl)ether	1.000	1.170	-17.0	124	0.00
9		2-Methylphenol	1.000	1.073	-7.3	106	0.00
10		3+4-Methylphenols	1.000	1.101	-10.1	111	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	105	0.00
12	S	Nitrobenzene-d5	1.000	0.992	0.8	108	0.00
13		Nitrobenzene	1.000	0.942	5.8	101	0.00
14		2-Nitrophenol	1.000	1.039	-3.9	109	0.00
15		2,4-Dimethylphenol	1.000	1.049	-4.9	108	0.00
16	C	2,4-Dichlorophenol	1.000	1.044	-4.4	113	0.00
17	C	Hexachlorobutadiene	1.000	1.083	-8.3	111	-0.03
18		4-Chloro-3-methylphenol	1.000	1.074	-7.4	114	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	109	-0.02
20	S	2,4,6-Tribromophenol	1.000	0.904	9.6	104	0.00
21		2,4,6-Trichlorophenol	1.000	0.930	7.0	103	0.00
22		2,4,5-Trichlorophenol	1.000	1.046	-4.6	107	0.00
23	S	2-Fluorobiphenyl	1.000	1.024	-2.4	110	0.00
24	P	2,4-Dinitrophenol	1.000	0.000	100.0#	0	-14.40#
25		4-Nitrophenol	1.000	0.929	7.1	102	0.00
26	I	Phenanthrene-d10	5.000	5.000	0.0	118	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.933	6.7	116	0.00
28		Hexachlorobenzene	1.000	0.853	14.7	100	0.00
29	C	Pentachlorophenol	1.000	0.843	15.7	88	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	107	0.00
31		Benzidine	1.000	1.795	-79.5#	1370	-0.02
32	S	Terphenyl-d14	1.000	1.033	-3.3	108	0.00
33		3,3'-Dichlorobenzidine	1.000	1.001	-0.1	144	0.00
34	I	Perylene-d12	5.000	5.000	0.0	107	0.00

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