Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040315\
Data File : BM000839.D
Acq On : 03 Apr 2015 13:06
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Apr 03 18:52:16 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	99	0.00
2		1,4-Dioxane	1.000	1.081	-8.1	103	0.00
3		n-Nitrosodimethylamine	1.000	1.020	-2.0	100	0.00
4	S	2-Fluorophenol	1.000	1.052	-5.2	103	0.00
5	S	Phenol-d6	1.000	1.007	-0.7	105	0.00
6		2-Chlorophenol	1.000	1.038	-3.8	107	0.00
7	C	Phenol	1.000	1.008	-0.8	101	0.00
8		bis(2-Chloroethyl)ether	1.000	1.013	-1.3	108	0.00
9		2-Methylphenol	1.000	0.995	0.5	101	0.00
10		3+4-Methylphenols	1.000	0.975	2.5	100	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	101	0.00
12	S	Nitrobenzene-d5	1.000	0.970	3.0	102	0.00
13		Nitrobenzene	1.000	0.938	6.2	98	0.00
14		2-Nitrophenol	1.000	0.977	2.3	98	0.00
15		2,4-Dimethylphenol	1.000	1.021	-2.1	101	0.00
16	C	2,4-Dichlorophenol	1.000	0.984	1.6	101	0.00
17	C	Hexachlorobutadiene	1.000	1.051	-5.1	104	-0.03
18		4-Chloro-3-methylphenol	1.000	1.008	-0.8	103	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
20	S	2,4,6-Tribromophenol	1.000	0.867	13.3	94	0.00
21		2,4,6-Trichlorophenol	1.000	0.854	14.6	89	0.00
22		2,4,5-Trichlorophenol	1.000	0.994	0.6	96	0.00
23	S	2-Fluorobiphenyl	1.000	1.050	-5.0	107	0.00
24	P	2,4-Dinitrophenol	1.000	0.213	78.7#	48	0.00
25		4-Nitrophenol	1.000	0.827	17.3	82	0.00
26	I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.790	21.0	81	0.00
28		Hexachlorobenzene	1.000	0.928	7.2	101	0.00
29	C	Pentachlorophenol	1.000	0.690	31.0#	58	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	102	0.00
31		Benzidine	1.000	1.409	-40.9#	898	-0.02
32	S	Terphenyl-d14	1.000	1.036	-3.6	104	0.00
33		3,3'-Dichlorobenzidine	1.000	0.909	9.1	123	0.00
34	I	Perylene-d12	5.000	5.000	0.0	103	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1