Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000828.D
Acq On : 03 Apr 2015 04:08
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Apr 03 18:39:34 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	92	0.00
2		1,4-Dioxane	1.000	1.084	-8.4	96	0.00
3		n-Nitrosodimethylamine	1.000	1.049	-4.9	96	0.00
4	S	2-Fluorophenol	1.000	1.078	-7.8	98	0.00
5	S	Phenol-d6	1.000	1.016	-1.6	99	0.00
6		2-Chlorophenol	1.000	1.036	-3.6	100	0.00
7	C	Phenol	1.000	1.023	-2.3	96	0.00
8		bis(2-Chloroethyl)ether	1.000	1.019	-1.9	101	0.00
9		2-Methylphenol	1.000	1.007	-0.7	95	0.00
10		3+4-Methylphenols	1.000	0.988	1.2	95	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	98	0.00
12	S	Nitrobenzene-d5	1.000	0.956	4.4	97	0.00
13		Nitrobenzene	1.000	0.942	5.8	94	0.00
14		2-Nitrophenol	1.000	0.981	1.9	95	0.00
15		2,4-Dimethylphenol	1.000	1.001	-0.1	96	0.00
16	C	2,4-Dichlorophenol	1.000	0.963	3.7	95	0.00
17	C	Hexachlorobutadiene	1.000	1.036	-3.6	99	-0.03
18		4-Chloro-3-methylphenol	1.000	1.003	-0.3	99	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
20	S	2,4,6-Tribromophenol	1.000	0.849	15.1	89	-0.02
21		2,4,6-Trichlorophenol	1.000	0.877	12.3	88	0.00
22		2,4,5-Trichlorophenol	1.000	1.006	-0.6	94	0.00
23	S	2-Fluorobiphenyl	1.000	1.010	-1.0	100	0.00
24	P	2,4-Dinitrophenol	1.000	0.257	74.3#	56	0.01
25		4-Nitrophenol	1.000	0.884	11.6	87	0.01
26	I	Phenanthrene-d10	5.000	5.000	0.0	121	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.817	18.3	96	0.00
28		Hexachlorobenzene	1.000	0.721	27.9#	85	-0.02
29	C	Pentachlorophenol	1.000	0.716	28.4#	69	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
31		Benzidine	1.000	1.200	-20.0	696	-0.03
32	S	Terphenyl-d14	1.000	1.026	-2.6	102	0.00
33		3,3'-Dichlorobenzidine	1.000	0.900	10.0	120	0.00
34	I	Perylene-d12	5.000	5.000	0.0	102	0.00

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