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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	95	0.00
2		1,4-Dioxane	0.756	0.779	-3.0	107	0.00
3		n-Nitrosodimethylamine	0.635	0.670	-5.5	100	0.00
4	S	2-Fluorophenol	1.377	1.511	-9.7	104	0.00
5	S	Phenol-d6	1.921	2.222	-15.7	109	0.00
6		2-Chlorophenol	1.365	1.588	-16.3	111	0.00
7	C	Phenol	2.063	2.125	-3.0	100	0.00
8		bis(2-Chloroethyl)ether	1.435	1.554	-8.3	109	0.00
9		2-Methylphenol	1.472	1.592	-8.2	103	0.00
10		3+4-Methylphenols	1.558	1.827	-17.3	103	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
12	S	Nitrobenzene-d5	0.233	0.230	1.3	103	0.00
13		Nitrobenzene	0.272	0.262	3.7	100	0.00
14		2-Nitrophenol	0.142	0.136	4.2	103	0.00
15		2,4-Dimethylphenol	0.283	0.294	-3.9	102	0.00
16	C	2,4-Dichlorophenol	0.251	0.263	-4.8	103	0.00
17	C	Hexachlorobutadiene	0.193	0.201	-4.1	103	-0.03
18		4-Chloro-3-methylphenol	0.251	0.259	-3.2	105	0.00
19	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
20	S	2,4,6-Tribromophenol	0.149	0.155	-4.0	99	0.00
21		2,4,6-Trichlorophenol	0.270	0.294	-8.9	94	0.00
22		2,4,5-Trichlorophenol	0.331	0.385	-16.3	102	0.00
23	S	2-Fluorobiphenyl	1.540	1.557	-1.1	103	0.00
24	P	2,4-Dinitrophenol	0.049	0.014#	71.4#	65	0.00
25		4-Nitrophenol	0.098	0.087	11.2	95	0.00
26	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
27		4,6-Dinitro-2-methylphenol	0.044	0.036	18.2	82	0.03
28		Hexachlorobenzene	0.282	0.272	3.5	100	0.00
29	C	Pentachlorophenol	0.049	0.037	24.5#	75	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
31		Benzidine	0.201	0.293	-45.8#	1187#	-0.02
32	S	Terphenyl-d14	0.602	0.614	-2.0	103	0.00
33		3,3'-Dichlorobenzidine	0.336	0.375	-11.6	138	0.00
34	I	Perylene-d12	1.000	1.000	0.0	104	0.00

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