

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