

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	AvqRF	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