

Data Path : Z:\HPCHEM1\BNA M\Data\BM040715\
 Data File : BM000900.D
 Acq On : 08 Apr 2015 04:20
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Apr 08 06:04:23 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 07 15:59:39 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	109	-0.04
2	1,4-Dioxane	0.864	0.809	6.4	104	0.00
3	n-Nitrosodimethylamine	0.746	0.720	3.5	111	0.00
4 S	2-Fluorophenol	1.678	1.655	1.4	111	0.00
5 S	Phenol-d6	2.645	2.577	2.6	113	0.00
6	2-Chlorophenol	2.076	2.182	-5.1	122	0.00
7 C	Phenol	2.091	1.794	14.2	99	0.00
8	bis(2-Chloroethyl)ether	2.038	2.179	-6.9	125	0.00
9	2-Methylphenol	1.533	1.305	14.9	95	0.00
10	3+4-Methylphenols	1.999	1.851	7.4	108	0.00
11 I	Naphthalene-d8	1.000	1.000	0.0	108	-0.03
12 S	Nitrobenzene-d5	0.249	0.241	3.2	112	-0.04
13	Nitrobenzene	0.251	0.222	11.6	99	0.00
14	2-Nitrophenol	0.158	0.108	31.6#	84	0.00
15	2,4-Dimethylphenol	0.287	0.265	7.7	101	0.00
16 C	2,4-Dichlorophenol	0.281	0.287	-2.1	116	0.00
17 C	Hexachlorobutadiene	0.228	0.249	-9.2	117	0.00
18	4-Chloro-3-methylphenol	0.274	0.265	3.3	112	0.00
19 I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
20 S	2,4,6-Tribromophenol	0.145	0.132	9.0	103	0.02
21	2,4,6-Trichlorophenol	0.256	0.297	-16.0	134	0.00
22	2,4,5-Trichlorophenol	0.350	0.373	-6.6	127	0.00
23 S	2-Fluorobiphenyl	1.541	1.535	0.4	109	0.00
24 P	2,4-Dinitrophenol	0.000	0.000#	0.0	0#	-14.40#
25	4-Nitrophenol	0.061	0.064	-4.9	95	-0.01
26 I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
27	4,6-Dinitro-2-methylphenol	0.049	0.034	30.6#	94	0.00
28	Hexachlorobenzene	0.222	0.192	13.5	85	0.00
29 C	Pentachlorophenol	0.052	0.029	44.2#	166#	0.00
30 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
31	Benzidine	0.222	0.213	4.1	102	0.00
32 S	Terphenyl-d14	0.617	0.618	-0.2	106	0.00
33	3,3'-Dichlorobenzidine	0.330	0.312	5.5	100	0.00
34 I	Perylene-d12	1.000	1.000	0.0	105	0.00

(#) = Out of Range

SPCC's out = 1 CCC's out = 1