

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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
2	1,4-Dioxane	0.756	0.725	4.1	96	0.00
3	n-Nitrosodimethylamine	0.635	0.667	-5.0	96	0.00
4 S	2-Fluorophenol	1.377	1.484	-7.8	98	0.00
5 S	Phenol-d6	1.921	2.087	-8.6	99	0.00
6	2-Chlorophenol	1.365	1.478	-8.3	100	0.00
7 C	Phenol	2.063	2.111	-2.3	96	0.00
8	bis(2-Chloroethyl)ether	1.435	1.495	-4.2	101	0.00
9	2-Methylphenol	1.472	1.537	-4.4	95	0.00
10	3+4-Methylphenols	1.558	1.752	-12.5	95	0.00
11 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
12 S	Nitrobenzene-d5	0.233	0.223	4.3	97	0.00
13	Nitrobenzene	0.272	0.257	5.5	94	0.00
14	2-Nitrophenol	0.142	0.129	9.2	95	0.00
15	2,4-Dimethylphenol	0.283	0.284	-0.4	96	0.00
16 C	2,4-Dichlorophenol	0.251	0.250	0.4	95	0.00
17 C	Hexachlorobutadiene	0.193	0.200	-3.6	99	-0.03
18	4-Chloro-3-methylphenol	0.251	0.252	-0.4	99	0.00
19 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
20 S	2,4,6-Tribromophenol	0.149	0.144	3.4	89	-0.02
21	2,4,6-Trichlorophenol	0.270	0.286	-5.9	88	0.00
22	2,4,5-Trichlorophenol	0.331	0.368	-11.2	94	0.00
23 S	2-Fluorobiphenyl	1.540	1.555	-1.0	100	0.00
24 P	2,4-Dinitrophenol	0.049	0.013#	73.5#	56	0.01
25	4-Nitrophenol	0.098	0.082	16.3	87	0.01
26 I	Phenanthrene-d10	1.000	1.000	0.0	121	0.00
27	4,6-Dinitro-2-methylphenol	0.044	0.037	15.9	96	0.00
28	Hexachlorobenzene	0.282	0.203	28.0#	85	-0.02
29 C	Pentachlorophenol	0.049	0.030	38.8#	69	0.00
30 I	Chrysene-d12	1.000	1.000	0.0	101	0.00
31	Benzidine	0.201	0.175	12.9	696#	-0.03
32 S	Terphenyl-d14	0.602	0.617	-2.5	102	0.00
33	3,3'-Dichlorobenzidine	0.336	0.333	0.9	120	0.00
34 I	Perylene-d12	1.000	1.000	0.0	102	0.00

(#) = Out of Range

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