

Data Path : Z:\HPCHEM1\BNA M\Data\BM040715\
 Data File : BM000887.D
 Acq On : 07 Apr 2015 19:25
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Apr 08 01:55:52 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	108	0.00
2	1,4-Dioxane	1.000	1.047	-4.7	107	0.00
3	n-Nitrosodimethylamine	1.000	0.974	2.6	116	-0.02
4 S	2-Fluorophenol	1.000	1.001	-0.1	112	0.00
5 S	Phenol-d6	1.000	0.962	3.8	116	0.00
6	2-Chlorophenol	1.000	0.963	3.7	118	0.00
7 C	Phenol	1.000	0.889	11.1	102	0.00
8	bis(2-Chloroethyl)ether	1.000	0.967	3.3	122	0.00
9	2-Methylphenol	1.000	0.952	4.8	105	0.00
10	3+4-Methylphenols	1.000	0.941	5.9	113	0.00
11 I	Naphthalene-d8	5.000	5.000	0.0	112	0.00
12 S	Nitrobenzene-d5	1.000	0.939	6.1	113	-0.04
13	Nitrobenzene	1.000	0.913	8.7	105	0.00
14	2-Nitrophenol	1.000	0.989	1.1	113	0.00
15	2,4-Dimethylphenol	1.000	0.968	3.2	110	0.00
16 C	2,4-Dichlorophenol	1.000	0.934	6.6	115	0.00
17 C	Hexachlorobutadiene	1.000	1.021	-2.1	113	0.00
18	4-Chloro-3-methylphenol	1.000	0.951	4.9	115	0.00
19 I	Acenaphthene-d10	5.000	5.000	0.0	113	0.00
20 S	2,4,6-Tribromophenol	1.000	0.873	12.7	102	0.02
21	2,4,6-Trichlorophenol	1.000	0.872	12.8	124	0.00
22	2,4,5-Trichlorophenol	1.000	0.902	9.8	123	0.00
23 S	2-Fluorobiphenyl	1.000	1.001	-0.1	114	0.00
24 P	2,4-Dinitrophenol	1.000	0.000	100.0#	0	-14.40#
25	4-Nitrophenol	1.000	0.933	6.7	107	-0.01
26 I	Phenanthrene-d10	5.000	5.000	0.0	125	0.00
27	4,6-Dinitro-2-methylphenol	1.000	0.783	21.7	96	0.00
28	Hexachlorobenzene	1.000	0.879	12.1	87	0.02
29 C	Pentachlorophenol	1.000	0.907	9.3	145	0.00
30 I	Chrysene-d12	5.000	5.000	0.0	104	0.00
31	Benzidine	1.000	0.998	0.2	108	0.00
32 S	Terphenyl-d14	1.000	1.002	-0.2	104	0.00
33	3,3'-Dichlorobenzidine	1.000	1.014	-1.4	110	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