

Data Path : Z:\HPCHEM1\BNA M\Data\BM050615\
 Data File : BM001233.D
 Acq On : 07 May 2015 00:52
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM050615

Quant Time: May 07 07:18:55 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM050615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 07 07:00:24 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	126	0.00
2	1,4-Dioxane	1.000	1.024	-2.4	130	0.00
3 S	2-Fluorophenol	1.000	1.048	-4.8	130	0.00
4 S	Phenol-d6	1.000	1.025	-2.5	132	0.00
5	2-Chlorophenol	1.000	1.017	-1.7	128	0.00
6 C	Phenol	1.000	0.992	0.8	123	0.00
7	2-Methylphenol	1.000	1.004	-0.4	125	0.00
8	3+4-Methylphenols	1.000	1.031	-3.1	136	-0.03
9 I	Naphthalene-d8	5.000	5.000	0.0	130	0.00
10 S	Nitrobenzene-d5	1.000	0.999	0.1	136	0.00
11 C	2-Nitrophenol	1.000	0.953	4.7	130	0.00
12	2,4-Dimethylphenol	1.000	1.006	-0.6	129	0.00
13 C	2,4-Dichlorophenol	1.000	1.008	-0.8	129	0.00
14 C	4-Chloro-3-methylphenol	1.000	0.960	4.0	129	0.00
15	2-Methylnaphthalene	1.000	1.030	-3.0	130	0.00
16 I	Acenaphthene-d10	5.000	5.000	0.0	131	0.00
17 S	2,4,6-Tribromophenol	1.000	0.946	5.4	131	0.00
18 C	2,4,6-Trichlorophenol	1.000	0.972	2.8	131	0.00
19	2,4,5-Trichlorophenol	1.000	0.960	4.0	131	0.00
20 S	2-Fluorobiphenyl	1.000	1.031	-3.1	131	0.00
21 P	2,4-Dinitrophenol	1.000	0.953	4.7	133	0.00
22 P	4-Nitrophenol	1.000	1.095	-9.5	145	-0.01
23 I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
24	4,6-Dinitro-2-methylphenol	1.000	0.962	3.8	131	0.00
25	Hexachlorobenzene	1.000	1.033	-3.3	132	0.00
26 C	Pentachlorophenol	1.000	1.023	-2.3	126	0.00
27 I	Chrysene-d12	5.000	5.000	0.0	136	0.00
28 S	Terphenyl-d14	1.000	1.004	-0.4	133	0.00
29 I	Perylene-d12	5.000	5.000	0.0	138	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 0