

Data Path : Z:\HPCHEM1\BNA_E\Data\BE120914\
 Data File : BE089187.D
 Acq On : 9 Dec 2014 11:02
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Dec 10 00:35:11 2014
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIM-ACID-BE120814.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 09 19:38:28 2014
 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	112	0.00
2 S	2-Fluorophenol	1.000	1.043	-4.3	107	0.00
3 S	Phenol-d6	1.000	1.050	-5.0	111	0.00
4	2-Chlorophenol	1.000	1.049	-4.9	112	0.00
5 C	Phenol	1.000	1.063	-6.3	112	0.00
6	2-Methylphenol	1.000	1.047	-4.7	108	0.00
7	3+4-Methylphenols	1.000	1.066	-6.6	113	0.01
8 I	Naphthalene-d8	5.000	5.000	0.0	112	0.00
9 C	2-Nitrophenol	1.000	1.013	-1.3	109	0.00
10	2,4-Dimethylphenol	1.000	1.031	-3.1	112	0.00
11 C	2,4-Dichlorophenol	1.000	1.034	-3.4	114	0.00
12 C	4-Chloro-3-methylphenol	1.000	1.034	-3.4	113	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	113	0.00
14 S	2,4,6-Tribromophenol	1.000	1.038	-3.8	110	0.00
15 C	2,4,6-Trichlorophenol	1.000	1.032	-3.2	114	0.00
16	2,4,5-Trichlorophenol	1.000	1.037	-3.7	111	0.00
17 P	2,4-Dinitrophenol	1.000	0.975	2.5	118	0.01
18 P	4-Nitrophenol	1.000	1.084	-8.4	113	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	110	0.00
20	4,6-Dinitro-2-methylphenol	1.000	0.942	5.8	115	0.00
21 C	Pentachlorophenol	1.000	0.935	6.5	110	0.00
22 I	Chrysene-d12	5.000	5.000	0.0	105	0.00
23 I	Perylene-d12	5.000	5.000	0.0	104	0.00

(#) = Out of Range

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