

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101415\
 Data File : BM002904.D
 Acq On : 14 Oct 2015 19:28
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 15 01:15:47 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 01:02:42 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	91	0.00
2 S	2-Fluorophenol	1.000	0.919	8.1	88	0.02
3 S	Phenol-d6	1.000	0.865	13.5	83	0.02
4	2-Chlorophenol	1.000	0.890	11.0	82	0.00
5 C	Phenol	1.000	0.903	9.7	83	0.00
6	2-Methylphenol	1.000	0.902	9.8	85	0.00
7	3+4-Methylphenols	1.000	0.884	11.6	82	0.00
8 I	Naphthalene-d8	5.000	5.000	0.0	90	0.00
9 S	Nitrobenzene-d5	1.000	0.926	7.4	83	0.00
10 C	2-Nitrophenol	1.000	0.826	17.4	77	0.00
11	2,4-Dimethylphenol	1.000	0.883	11.7	81	0.00
12 C	2,4-Dichlorophenol	1.000	0.927	7.3	81	0.00
13 C	4-Chloro-3-methylphenol	1.000	0.870	13.0	81	0.00
14 I	Acenaphthene-d10	5.000	5.000	0.0	89	0.00
15 S	2,4,6-Tribromophenol	1.000	0.852	14.8	79	0.00
16 C	2,4,6-Trichlorophenol	1.000	0.866	13.4	79	0.00
17	2,4,5-Trichlorophenol	1.000	0.885	11.5	80	0.00
18 S	2-Fluorobiphenyl	1.000	0.929	7.1	84	0.01
19 P	2,4-Dinitrophenol	1.000	0.902	9.8	75	0.02
20 P	4-Nitrophenol	1.000	0.893	10.7	75	0.00
21 I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
22	4,6-Dinitro-2-methylphenol	1.000	0.848	15.2	73	0.00
23 C	Pentachlorophenol	1.000	1.025	-2.5	83	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	93	0.00
25 S	Terphenyl-d14	1.000	0.927	7.3	89	0.00
26 I	Perylene-d12	5.000	5.000	0.0	92	0.00

(#) = Out of Range

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