

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
 Data File : BM002899.D
 Acq On : 13 Oct 2015 22:48
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM101315

Quant Time: Oct 14 17:42:48 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 14 17:36:42 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	87	0.00
2 S	2-Fluorophenol	0.963	0.823	14.5	78	0.02
3 S	Phenol-d6	1.373	1.200	12.6	80	0.02
4	2-Chlorophenol	1.272	1.149	9.7	80	0.00
5 C	Phenol	1.349	1.229	8.9	80	0.00
6	2-Methylphenol	1.032	0.917	11.1	80	0.00
7	3+4-Methylphenols	1.277	1.132	11.4	78	0.01
8 I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
9 S	Nitrobenzene-d5	0.276	0.260	5.8	81	0.00
10 C	2-Nitrophenol	0.126	0.109	13.5	77	0.00
11	2,4-Dimethylphenol	0.276	0.250	9.4	79	0.00
12 C	2,4-Dichlorophenol	0.215	0.198	7.9	76	0.02
13 C	4-Chloro-3-methylphenol	0.269	0.237	11.9	78	0.00
14 I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
15 S	2,4,6-Tribromophenol	0.183	0.157	14.2	77	0.00
16 C	2,4,6-Trichlorophenol	0.326	0.277	15.0	75	0.00
17	2,4,5-Trichlorophenol	0.347	0.305	12.1	77	0.00
18 S	2-Fluorobiphenyl	1.636	1.516	7.3	81	0.01
19 P	2,4-Dinitrophenol	0.027	0.007#	74.1#	55	0.02
20 P	4-Nitrophenol	0.132	0.089	32.6#	69	0.00
21 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
22	4,6-Dinitro-2-methylphenol	0.025	0.025	0.0	73	0.00
23 C	Pentachlorophenol	0.038	0.023	39.5#	69	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	93	0.00
25 S	Terphenyl-d14	0.760	0.683	10.1	86	0.00
26 I	Perylene-d12	1.000	1.000	0.0	91	0.00

(#) = Out of Range

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