

Data Path : Z:\HPCHEM1\BNA M\Data\BM051615\
 Data File : BM001320.D
 Acq On : 15 May 2015 08:04
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 15 08:52:04 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM051315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 05:24:15 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	95	-0.02
2	1,4-Dioxane	1.000	0.905	9.5	90	0.00
3	n-Nitrosodimethylamine	1.000	0.883	11.7	96	0.00
4 S	2-Fluorophenol	1.000	0.932	6.8	93	-0.02
5 S	Phenol-d6	1.000	0.922	7.8	95	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
7 S	Nitrobenzene-d5	1.000	0.860	14.0	92	0.00
8	Nitrobenzene	1.000	0.849	15.1	93	0.00
9	Naphthalene	1.000	0.965	3.5	95	0.00
10	Hexachlorobutadiene	1.000	0.957	4.3	97	0.00
11	2-Methylnaphthalene	1.000	0.935	6.5	94	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
13 S	2,4,6-Tribromophenol	1.000	0.702	29.8#	79	-0.03
14 S	2-Fluorobiphenyl	1.000	0.957	4.3	94	0.00
15	Acenaphthylene	1.000	0.932	6.8	93	0.00
16	Acenaphthene	1.000	0.936	6.4	94	0.00
17	Fluorene	1.000	0.943	5.7	99	-0.03
18 I	Phenanthrene-d10	5.000	5.000	0.0	124	0.00
19	Hexachlorobenzene	1.000	0.646	35.4#	86	0.00
20	Pentachlorophenol	1.000	0.868	13.2	117	0.00
21	Phenanthrene	1.000	1.062	-6.2	143	0.00
22	Anthracene	1.000	1.007	-0.7	136	0.00
23	Fluoranthene	1.000	0.685	31.5#	94	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	97	-0.01
25	Pyrene	1.000	0.898	10.2	96	0.00
26 S	Terphenyl-d14	1.000	0.897	10.3	93	-0.01
27	Benzo(a)anthracene	1.000	0.909	9.1	100	-0.01
28	Chrysene	1.000	0.939	6.1	97	-0.01
29	Indeno(1,2,3-cd)pyrene	1.000	1.027	-2.7	107	-0.01
30 I	Perylene-d12	5.000	5.000	0.0	102	-0.02
31	Benzo(b)fluoranthene	1.000	0.903	9.7	101	-0.01
32	Benzo(k)fluoranthene	1.000	0.925	7.5	101	-0.01
33 C	Benzo(a)pyrene	1.000	0.923	7.7	104	-0.01
34	Dibenzo(a,h)anthracene	1.000	0.981	1.9	107	-0.02
35	Benzo(q,h,i)perylene	1.000	0.983	1.7	108	-0.02

(#) = Out of Range

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