

Data Path : Z:\HPCHEM1\BNA_M\Data\BM040115\
 Data File : BM000790.D
 Acq On : 01 Apr 2015 17:54
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Apr 02 01:57:22 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM033115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 01 04:17:30 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	94	0.00
2	1,4-Dioxane	1.000	0.000	100.0#	0	-2.47#
3	n-Nitrosodimethylamine	1.000	0.961	3.9	91	0.00
4 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
5 S	Nitrobenzene-d5	1.000	0.992	0.8	95	0.00
6	Naphthalene	1.000	1.026	-2.6	93	0.00
7	2-Methylnaphthalene	1.000	1.025	-2.5	93	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
9 S	2-Fluorobiphenyl	1.000	1.033	-3.3	93	0.00
10	Acenaphthylene	1.000	1.021	-2.1	95	0.00
11	Acenaphthene	1.000	1.019	-1.9	96	0.00
12	Fluorene	1.000	1.054	-5.4	97	0.00
13 I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
14	Hexachlorobenzene	1.000	0.980	2.0	96	0.00
15	Pentachlorophenol	1.000	0.792	20.8	72	0.00
16	Phenanthrene	1.000	1.004	-0.4	100	0.00
17	Anthracene	1.000	0.973	2.7	97	0.00
18	Fluoranthene	1.000	1.000	0.0	101	0.00
19 I	Chrysene-d12	5.000	5.000	0.0	103	0.00
20	Pyrene	1.000	1.001	-0.1	102	0.00
21 S	Terphenyl-d14	1.000	1.019	-1.9	102	0.00
22	Benzo(a)anthracene	1.000	0.997	0.3	106	-0.01
23	Chrysene	1.000	1.013	-1.3	102	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	1.019	-1.9	106	-0.01
25 I	Perylene-d12	5.000	5.000	0.0	107	0.00
26	Benzo(b)fluoranthene	1.000	1.067	-6.7	112	-0.01
27	Benzo(k)fluoranthene	1.000	0.943	5.7	98	0.00
28 C	Benzo(a)pyrene	1.000	1.015	-1.5	108	-0.01
29	Dibenzo(a,h)anthracene	1.000	1.001	-0.1	105	-0.01
30	Benzo(g,h,i)perylene	1.000	1.002	-0.2	105	-0.01

(#) = Out of Range

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