

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM033115\
 Data File : BM000775.D
 Acq On : 01 Apr 2015 04:22
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Apr 01 15:09:42 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	93	0.00
2	1,4-Dioxane	1.000	0.000	100.0#	0	-2.47#
3	n-Nitrosodimethylamine	1.000	0.971	2.9	92	0.00
4 I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
5 S	Nitrobenzene-d5	1.000	0.957	4.3	91	0.00
6	Naphthalene	1.000	1.028	-2.8	93	0.00
7	2-Methylnaphthalene	1.000	1.019	-1.9	93	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
9 S	2-Fluorobiphenyl	1.000	1.032	-3.2	93	0.00
10	Acenaphthylene	1.000	0.986	1.4	92	0.00
11	Acenaphthene	1.000	0.991	0.9	93	0.00
12	Fluorene	1.000	1.025	-2.5	94	0.00
13 I	Phenanthrene-d10	5.000	5.000	0.0	97	0.00
14	Hexachlorobenzene	1.000	1.007	-0.7	95	0.00
15	Pentachlorophenol	1.000	0.759	24.1	64	0.00
16	Phenanthrene	1.000	1.004	-0.4	96	0.00
17	Anthracene	1.000	0.955	4.5	92	0.00
18	Fluoranthene	1.000	0.970	3.0	94	0.00
19 I	Chrysene-d12	5.000	5.000	0.0	96	0.00
20	Pyrene	1.000	1.006	-0.6	95	0.00
21 S	Terphenyl-d14	1.000	1.024	-2.4	95	0.00
22	Benzo(a)anthracene	1.000	0.972	2.8	96	-0.01
23	Chrysene	1.000	1.032	-3.2	97	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	1.027	-2.7	99	-0.01
25 I	Perylene-d12	5.000	5.000	0.0	98	0.00
26	Benzo(b)fluoranthene	1.000	1.012	-1.2	98	0.00
27	Benzo(k)fluoranthene	1.000	1.018	-1.8	97	0.00
28 C	Benzo(a)pyrene	1.000	0.991	0.9	97	-0.01
29	Dibenzo(a,h)anthracene	1.000	1.027	-2.7	99	0.00
30	Benzo(g,h,i)perylene	1.000	1.031	-3.1	99	0.00

(#) = Out of Range

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