

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM060215\
 Data File : BM001532.D
 Acq On : 03 Jun 2015 02:36
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jun 03 03:36:33 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM060215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 03 02:19:00 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	0.00
2	1,4-Dioxane	1.000	0.993	0.7	98	0.00
3	n-Nitrosodimethylamine	1.000	1.034	-3.4	96	0.00
4 S	2-Fluorophenol	1.000	1.018	-1.8	99	-0.02
5 S	Phenol-d6	1.000	1.009	-0.9	98	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
7 S	Nitrobenzene-d5	1.000	1.021	-2.1	99	0.00
8	Nitrobenzene	1.000	1.015	-1.5	98	0.00
9	Naphthalene	1.000	0.983	1.7	94	0.00
10	Hexachlorobutadiene	1.000	0.998	0.2	95	0.00
11	2-Methylnaphthalene	1.000	0.990	1.0	94	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	95	0.00
13 S	2,4,6-Tribromophenol	1.000	1.254	-25.4#	119	0.00
14 S	2-Fluorobiphenyl	1.000	0.983	1.7	94	0.00
15	Acenaphthylene	1.000	0.978	2.2	94	0.00
16	Acenaphthene	1.000	0.987	1.3	93	0.00
17	Fluorene	1.000	0.840	16.0	81	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
19	4-Bromophenyl-phenylether	1.000	0.949	5.1	97	0.00
20	Hexachlorobenzene	1.000	0.723	27.7#	75	0.00
21	Pentachlorophenol	1.000	0.808	19.2	82	0.00
22	Phenanthrene	1.000	1.158	-15.8	123	0.00
23	Anthracene	1.000	0.759	24.1	78	0.00
24	Fluoranthene	1.000	0.904	9.6	95	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	95	-0.01
26	Pyrene	1.000	0.991	0.9	96	0.00
27 S	Terphenyl-d14	1.000	0.983	1.7	93	0.00
28	Benzo(a)anthracene	1.000	1.011	-1.1	100	-0.01
29	Chrysene	1.000	0.982	1.8	96	-0.01
30	Bis(2-ethylhexyl)phthalate	1.000	0.955	4.5	104	-0.01
31	Indeno(1,2,3-cd)pyrene	1.000	1.009	-0.9	98	-0.02
32 I	Perylene-d12	5.000	5.000	0.0	98	0.00
33	Benzo(b)fluoranthene	1.000	0.979	2.1	93	0.00
34	Benzo(k)fluoranthene	1.000	1.009	-0.9	103	0.00
35 C	Benzo(a)pyrene	1.000	0.996	0.4	98	0.00
36	Dibenzo(a,h)anthracene	1.000	0.989	1.1	97	-0.02
37	Benzo(g,h,i)perylene	1.000	0.992	0.8	98	0.00

(#) = Out of Range

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