

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062215\
 Data File : BM001770.D
 Acq On : 22 Jun 2015 12:00
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jun 23 04:06:55 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 05:05:14 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	131	0.00
2	1,4-Dioxane	1.000	1.005	-0.5	127	0.00
3	n-Nitrosodimethylamine	1.000	1.058	-5.8	138	-0.02
4 S	2-Fluorophenol	1.000	1.024	-2.4	134	0.00
5 S	Phenol-d6	1.000	1.012	-1.2	135	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	129	0.00
7 S	Nitrobenzene-d5	1.000	1.024	-2.4	135	0.00
8	Nitrobenzene	1.000	1.028	-2.8	135	0.00
9	Naphthalene	1.000	0.990	1.0	129	0.00
10	Hexachlorobutadiene	1.000	1.003	-0.3	133	0.00
11	2-Methylnaphthalene	1.000	0.990	1.0	130	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	128	0.00
13 S	2,4,6-Tribromophenol	1.000	1.052	-5.2	138	0.00
14 S	2-Fluorobiphenyl	1.000	1.002	-0.2	128	0.00
15	Acenaphthylene	1.000	0.994	0.6	129	0.00
16	Acenaphthene	1.000	0.986	1.4	128	0.00
17	Fluorene	1.000	1.030	-3.0	118	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	114	0.00
19	4-Bromophenyl-phenylether	1.000	1.049	-4.9	132	0.00
20	Hexachlorobenzene	1.000	0.942	5.8	115	0.00
21	Pentachlorophenol	1.000	1.233	-23.3	151	0.00
22	Phenanthrene	1.000	0.964	3.6	116	0.00
23	Anthracene	1.000	0.996	0.4	113	0.00
24	Fluoranthene	1.000	1.038	-3.8	123	-0.01
25 I	Chrysene-d12	5.000	5.000	0.0	119	0.00
26	Pyrene	1.000	0.984	1.6	124	0.00
27 S	Terphenyl-d14	1.000	0.998	0.2	124	0.00
28	Benzo(a)anthracene	1.000	0.959	4.1	118	0.00
29	Chrysene	1.000	0.965	3.5	117	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.938	6.2	122	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.974	2.6	116	0.00
32 I	Perylene-d12	5.000	5.000	0.0	118	0.00
33	Benzo(b)fluoranthene	1.000	0.967	3.3	115	0.00
34	Benzo(k)fluoranthene	1.000	1.018	-1.8	123	0.00
35 C	Benzo(a)pyrene	1.000	1.001	-0.1	120	0.01
36	Dibenzo(a,h)anthracene	1.000	0.980	2.0	116	0.00
37	Benzo(g,h,i)perylene	1.000	0.980	2.0	117	0.00

(#) = Out of Range

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