

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050515\
 Data File : BM001223.D
 Acq On : 06 May 2015 12:39
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 06 13:24:56 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 12:46:35 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	105	0.00
2	1,4-Dioxane	0.522	0.504	3.4	89	0.00
3	n-Nitrosodimethylamine	0.607	0.554	8.7	106	0.00
4 S	2-Fluorophenol	1.045	1.054	-0.9	104	0.00
5 S	Phenol-d6	1.329	1.334	-0.4	102	-0.02
6 I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
7 S	Nitrobenzene-d5	0.283	0.287	-1.4	103	0.00
8	Nitrobenzene	0.300	0.303	-1.0	105	0.00
9	Naphthalene	1.089	1.122	-3.0	106	0.00
10	Hexachlorobutadiene	0.211	0.214	-1.4	107	0.00
11	2-Methylnaphthalene	0.765	0.791	-3.4	105	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
13 S	2,4,6-Tribromophenol	0.266	0.259	2.6	105	0.00
14 S	2-Fluorobiphenyl	1.529	1.588	-3.9	106	-0.01
15	Acenaphthylene	2.155	2.221	-3.1	105	0.00
16	Acenaphthene	1.355	1.388	-2.4	105	0.00
17	Fluorene	2.203	2.234	-1.4	99	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
19	Hexachlorobenzene	0.286	0.287	-0.3	105	0.00
20	Pentachlorophenol	0.107	0.082	23.4	94	0.00
21	Phenanthrene	2.093	2.104	-0.5	97	0.00
22	Anthracene	1.713	1.702	0.6	101	0.00
23	Fluoranthene	1.971	2.035	-3.2	103	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
25	Pyrene	1.598	1.593	0.3	104	0.00
26 S	Terphenyl-d14	0.716	0.713	0.4	103	0.00
27	Benzo(a)anthracene	1.426	1.442	-1.1	107	0.00
28	Chrysene	1.376	1.423	-3.4	106	0.00
29	Indeno(1,2,3-cd)pyrene	1.092	1.245	-14.0	117	0.00
30 I	Perylene-d12	1.000	1.000	0.0	113	0.00
31	Benzo(b)fluoranthene	1.599	1.522	4.8	111	0.00
32	Benzo(k)fluoranthene	1.469	1.573	-7.1	112	0.00
33 C	Benzo(a)pyrene	1.320	1.348	-2.1	113	0.00
34	Dibenzo(a,h)anthracene	1.074	1.132	-5.4	117	0.00
35	Benzo(g,h,i)perylene	1.114	1.187	-6.6	118	0.00

(#) = Out of Range

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