

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050415\
 Data File : BM001186.D
 Acq On : 04 May 2015 10:52
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 05 02:43:47 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 04 16:48:07 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	128	0.00
2	1,4-Dioxane	0.612	0.524	14.4	108	0.00
3	n-Nitrosodimethylamine	0.612	0.622	-1.6	124	0.00
4 S	2-Fluorophenol	0.972	1.014	-4.3	132	-0.01
5 S	Phenol-d6	1.198	1.293	-7.9	137	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	132	0.00
7 S	Nitrobenzene-d5	0.283	0.289	-2.1	135	0.00
8	Nitrobenzene	0.303	0.308	-1.7	127	0.00
9	Naphthalene	1.099	1.102	-0.3	130	0.00
10	Hexachlorobutadiene	0.212	0.211	0.5	132	0.00
11	2-Methylnaphthalene	0.754	0.780	-3.4	137	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
13 S	2,4,6-Tribromophenol	0.188	0.211	-12.2	164#	0.00
14 S	2-Fluorobiphenyl	1.654	1.693	-2.4	137	-0.01
15	Acenaphthylene	2.185	2.332	-6.7	143	-0.02
16	Acenaphthene	1.510	1.598	-5.8	146	0.00
17	Fluorene	1.751	2.036	-16.3	152#	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	144	0.00
19	Hexachlorobenzene	0.269	0.263	2.2	143	0.00
20	Pentachlorophenol	0.067	0.070	-4.5	157#	0.00
21	Phenanthrene	1.287	1.307	-1.6	147	0.00
22	Anthracene	1.064	1.075	-1.0	142	-0.02
23	Fluoranthene	1.303	1.331	-2.1	147	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	137	0.00
25	Pyrene	1.562	1.674	-7.2	147	0.00
26 S	Terphenyl-d14	0.689	0.730	-6.0	142	0.00
27	Benzo(a)anthracene	1.410	1.482	-5.1	145	0.00
28	Chrysene	1.384	1.404	-1.4	134	0.00
29	Indeno(1,2,3-cd)pyrene	1.262	1.110	12.0	116	0.00
30 I	Perylene-d12	1.000	1.000	0.0	124	-0.01
31	Benzo(b)fluoranthene	1.540	1.546	-0.4	115	0.00
32	Benzo(k)fluoranthene	1.442	1.602	-11.1	146	-0.01
33 C	Benzo(a)pyrene	1.299	1.370	-5.5	130	0.00
34	Dibenzo(a,h)anthracene	1.139	1.082	5.0	115	-0.01
35	Benzo(g,h,i)perylene	1.202	1.122	6.7	113	0.00

(#) = Out of Range

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