

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050415\
 Data File : BM001189.D
 Acq On : 04 May 2015 17:33
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 05 04:01:39 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	119	0.00
2	1,4-Dioxane	0.612	0.533	12.9	102	0.00
3	n-Nitrosodimethylamine	0.612	0.564	7.8	105	0.00
4 S	2-Fluorophenol	0.972	1.025	-5.5	124	-0.01
5 S	Phenol-d6	1.198	1.297	-8.3	128	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	123	0.00
7 S	Nitrobenzene-d5	0.283	0.290	-2.5	126	0.00
8	Nitrobenzene	0.303	0.323	-6.6	124	0.00
9	Naphthalene	1.099	1.109	-0.9	121	0.00
10	Hexachlorobutadiene	0.212	0.211	0.5	123	0.00
11	2-Methylnaphthalene	0.754	0.783	-3.8	128	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	130	0.00
13 S	2,4,6-Tribromophenol	0.188	0.211	-12.2	156#	0.00
14 S	2-Fluorobiphenyl	1.654	1.639	0.9	127	-0.01
15	Acenaphthylene	2.185	2.290	-4.8	135	-0.02
16	Acenaphthene	1.510	1.553	-2.8	135	0.00
17	Fluorene	1.751	1.977	-12.9	141	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	137	0.00
19	Hexachlorobenzene	0.269	0.261	3.0	135	0.00
20	Pentachlorophenol	0.067	0.071	-6.0	152#	0.00
21	Phenanthrene	1.287	1.285	0.2	137	0.00
22	Anthracene	1.064	1.091	-2.5	138	-0.02
23	Fluoranthene	1.303	1.334	-2.4	141	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	139	0.00
25	Pyrene	1.562	1.580	-1.2	141	0.00
26 S	Terphenyl-d14	0.689	0.705	-2.3	140	0.00
27	Benzo(a)anthracene	1.410	1.478	-4.8	147	0.00
28	Chrysene	1.384	1.400	-1.2	136	0.00
29	Indeno(1,2,3-cd)pyrene	1.262	1.265	-0.2	135	0.00
30 I	Perylene-d12	1.000	1.000	0.0	139	-0.01
31	Benzo(b)fluoranthene	1.540	1.545	-0.3	129	-0.01
32	Benzo(k)fluoranthene	1.442	1.532	-6.2	157#	-0.01
33 C	Benzo(a)pyrene	1.299	1.369	-5.4	145	0.00
34	Dibenzo(a,h)anthracene	1.139	1.123	1.4	134	-0.01
35	Benzo(g,h,i)perylene	1.202	1.162	3.3	131	0.00

(#) = Out of Range

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