

Data Path : Z:\HPCHEM1\BNA M\Data\BM050115\
 Data File : BM001172.D
 Acq On : 02 May 2015 17:37
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 04 17:40:18 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2	1,4-Dioxane	0.612	0.588	3.9	94	0.00
3	n-Nitrosodimethylamine	0.612	0.567	7.4	88	0.00
4 S	2-Fluorophenol	0.972	0.993	-2.2	100	-0.01
5 S	Phenol-d6	1.198	1.224	-2.2	101	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
7 S	Nitrobenzene-d5	0.283	0.281	0.7	99	0.00
8	Nitrobenzene	0.303	0.297	2.0	92	0.00
9	Naphthalene	1.099	1.114	-1.4	99	0.00
10	Hexachlorobutadiene	0.212	0.213	-0.5	100	0.00
11	2-Methylnaphthalene	0.754	0.778	-3.2	103	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
13 S	2,4,6-Tribromophenol	0.188	0.293	-55.9#	169#	0.00
14 S	2-Fluorobiphenyl	1.654	1.699	-2.7	102	0.00
15	Acenaphthylene	2.185	2.259	-3.4	103	-0.02
16	Acenaphthene	1.510	1.560	-3.3	106	0.00
17	Fluorene	1.751	1.933	-10.4	108	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
19	Hexachlorobenzene	0.269	0.263	2.2	104	0.00
20	Pentachlorophenol	0.067	0.062	7.5	101	0.00
21	Phenanthrene	1.287	1.310	-1.8	107	0.00
22	Anthracene	1.064	1.035	2.7	100	0.00
23	Fluoranthene	1.303	1.324	-1.6	107	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
25	Pyrene	1.562	1.588	-1.7	106	0.00
26 S	Terphenyl-d14	0.689	0.701	-1.7	104	0.00
27	Benzo(a)anthracene	1.410	1.424	-1.0	106	0.00
28	Chrysene	1.384	1.400	-1.2	102	0.00
29	Indeno(1,2,3-cd)pyrene	1.262	1.169	7.4	93	0.00
30 I	Perylene-d12	1.000	1.000	0.0	98	-0.01
31	Benzo(b)fluoranthene	1.540	1.514	1.7	89	-0.01
32	Benzo(k)fluoranthene	1.442	1.556	-7.9	113	-0.01
33 C	Benzo(a)pyrene	1.299	1.315	-1.2	99	0.00
34	Dibenzo(a,h)anthracene	1.139	1.097	3.7	93	-0.01
35	Benzo(q,h,i)perylene	1.202	1.144	4.8	91	0.00

(#) = Out of Range

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