

Data Path : Z:\HPCHEM1\BNA_M\Data\BM052015\
 Data File : BM001382.D
 Acq On : 22 May 2015 04:52
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 22 05:36:25 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 25 10:49:04 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	114	0.00
2	1,4-Dioxane	0.512	0.348	32.0#	75	0.00
3	n-Nitrosodimethylamine	0.564	0.708	-25.5#	135	-0.02
4 S	2-Fluorophenol	0.971	1.806	-86.0#	216#	0.00
5 S	Phenol-d6	1.208	2.156	-78.5#	206#	0.02
6 I	Naphthalene-d8	1.000	1.000	0.0	125	0.00
7 S	Nitrobenzene-d5	0.270	0.226	16.3	109	0.00
8	Nitrobenzene	0.282	0.237	16.0	110	0.00
9	Naphthalene	1.067	1.111	-4.1	131	0.00
10	Hexachlorobutadiene	0.206	0.196	4.9	120	0.00
11	2-Methylnaphthalene	0.758	0.816	-7.7	137	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	129	0.00
13 S	2,4,6-Tribromophenol	0.175	0.103	41.1#	74	0.00
14 S	2-Fluorobiphenyl	1.491	1.540	-3.3	134	0.00
15	Acenaphthylene	2.104	2.286	-8.7	145	0.00
16	Acenaphthene	1.321	1.386	-4.9	138	0.00
17	Fluorene	1.057	1.072	-1.4	134	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	131	0.00
19	4-Bromophenyl-phenylether	0.251	0.291	-15.9	145	0.00
20	Hexachlorobenzene	0.369	0.369	0.0	132	0.00
21	Pentachlorophenol	0.133	0.076	42.9#	94	0.00
22	Phenanthrene	1.758	1.761	-0.2	130	0.00
23	Anthracene	0.680	0.677	0.4	122	0.00
24	Fluoranthene	1.506	1.589	-5.5	137	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	101	-0.02
26	Pyrene	1.528	1.962	-28.4#	131	0.00
27 S	Terphenyl-d14	0.658	0.851	-29.3#	133	0.00
28	Benzo(a)anthracene	1.430	1.579	-10.4	117	0.00
29	Chrysene	1.369	1.398	-2.1	106	0.00
30	Bis(2-ethylhexyl)phthalate	0.705	1.284	-82.1#	205#	-0.01
31	Indeno(1,2,3-cd)pyrene	1.475	1.175	20.3	87	0.00
32 I	Perylene-d12	1.000	1.000	0.0	83	0.00
33	Benzo(b)fluoranthene	1.541	1.594	-3.4	89	0.00
34	Benzo(k)fluoranthene	1.357	1.594	-17.5	99	-0.01
35 C	Benzo(a)pyrene	1.311	1.438	-9.7	96	0.00
36	Dibenzo(a,h)anthracene	1.204	1.205	-0.1	88	0.00
37	Benzo(g,h,i)perylene	1.293	1.248	3.5	85	0.00

(#) = Out of Range

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