

Data Path : Z:\HPCHEM1\BNA M\DATA\BM052015\
 Data File : BM001363.D
 Acq On : 21 May 2015 16:20
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: May 21 17:05:38 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIMPAH-BM052015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 21 11:08:14 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	99	0.00
2	1,4-Dioxane	0.512	0.610	-19.1	113	0.00
3	n-Nitrosodimethylamine	0.564	0.619	-9.8	102	0.00
4 S	2-Fluorophenol	0.971	1.021	-5.1	105	0.00
5 S	Phenol-d6	1.208	1.305	-8.0	108	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
7 S	Nitrobenzene-d5	0.270	0.285	-5.6	109	0.00
8	Nitrobenzene	0.282	0.301	-6.7	111	0.00
9	Naphthalene	1.067	1.116	-4.6	104	0.00
10	Hexachlorobutadiene	0.206	0.216	-4.9	105	0.00
11	2-Methylnaphthalene	0.758	0.802	-5.8	106	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
13 S	2,4,6-Tribromophenol	0.175	0.178	-1.7	101	0.00
14 S	2-Fluorobiphenyl	1.491	1.554	-4.2	107	0.00
15	Acenaphthylene	2.104	2.201	-4.6	111	0.00
16	Acenaphthene	1.321	1.381	-4.5	109	0.00
17	Fluorene	1.057	1.099	-4.0	109	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
19	4-Bromophenyl-phenylether	0.251	0.247	1.6	105	0.00
20	Hexachlorobenzene	0.369	0.357	3.3	109	0.00
21	Pentachlorophenol	0.133	0.124	6.8	131	0.00
22	Phenanthrene	1.758	1.756	0.1	111	0.00
23	Anthracene	0.680	0.676	0.6	104	0.00
24	Fluoranthene	1.506	1.493	0.9	110	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	106	0.00
26	Pyrene	1.528	1.577	-3.2	111	0.00
27 S	Terphenyl-d14	0.658	0.677	-2.9	112	0.00
28	Benzo(a)anthracene	1.430	1.438	-0.6	113	0.00
29	Chrysene	1.369	1.402	-2.4	112	0.00
30	Bis(2-ethylhexyl)phthalate	0.705	0.797	-13.0	134	0.00
31	Indeno(1,2,3-cd)pyrene	1.475	1.497	-1.5	116	0.00
32 I	Perylene-d12	1.000	1.000	0.0	109	0.00
33	Benzo(b)fluoranthene	1.541	1.523	1.2	110	0.00
34	Benzo(k)fluoranthene	1.357	1.457	-7.4	118	0.00
35 C	Benzo(a)pyrene	1.311	1.359	-3.7	118	0.00
36	Dibenzo(a,h)anthracene	1.204	1.230	-2.2	116	0.01
37	Benzo(g,h,i)perylene	1.293	1.318	-1.9	117	0.00

(#) = Out of Range

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