

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061915\
 Data File : BM001752.D
 Acq On : 19 Jun 2015 00:22
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jun 19 05:17:37 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 19 05:05:14 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	105	0.00
2	1,4-Dioxane	0.570	0.578	-1.4	103	0.00
3	n-Nitrosodimethylamine	0.880	0.885	-0.6	105	-0.02
4 S	2-Fluorophenol	1.169	1.192	-2.0	107	0.00
5 S	Phenol-d6	1.505	1.491	0.9	106	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	105	-0.02
7 S	Nitrobenzene-d5	0.390	0.391	-0.3	107	0.00
8	Nitrobenzene	0.420	0.421	-0.2	107	0.00
9	Naphthalene	1.222	1.209	1.1	105	0.00
10	Hexachlorobutadiene	0.206	0.203	1.5	107	0.00
11	2-Methylnaphthalene	0.778	0.767	1.4	105	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	106	-0.01
13 S	2,4,6-Tribromophenol	0.156	0.158	-1.3	110	0.00
14 S	2-Fluorobiphenyl	1.787	1.772	0.8	105	0.00
15	Acenaphthylene	2.376	2.364	0.5	107	0.00
16	Acenaphthene	1.489	1.477	0.8	106	0.00
17	Fluorene	1.276	1.377	-7.9	102	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
19	4-Bromophenyl-phenylether	0.205	0.209	-2.0	109	0.00
20	Hexachlorobenzene	0.141	0.148	-5.0	109	0.00
21	Pentachlorophenol	0.063	0.070	-11.1	117	0.00
22	Phenanthrene	1.713	1.666	2.7	99	0.00
23	Anthracene	0.724	0.746	-3.0	99	0.00
24	Fluoranthene	1.080	1.148	-6.3	107	-0.01
25 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
26	Pyrene	1.741	1.698	2.5	108	-0.01
27 S	Terphenyl-d14	0.737	0.720	2.3	106	0.00
28	Benzo(a)anthracene	1.507	1.482	1.7	106	0.00
29	Chrysene	1.584	1.577	0.4	106	0.00
30	Bis(2-ethylhexyl)phthalate	0.869	0.837	3.7	110	0.00
31	Indeno(1,2,3-cd)pyrene	1.523	1.523	0.0	105	0.00
32 I	Perylene-d12	1.000	1.000	0.0	105	0.00
33	Benzo(b)fluoranthene	1.641	1.539	6.2	100	-0.01
34	Benzo(k)fluoranthene	1.469	1.554	-5.8	114	0.00
35 C	Benzo(a)pyrene	1.382	1.376	0.4	106	0.00
36	Dibenzo(a,h)anthracene	1.368	1.368	0.0	105	0.00
37	Benzo(g,h,i)perylene	1.439	1.422	1.2	105	0.00

(#) = Out of Range

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