

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM061215\
 Data File : BM001675.D
 Acq On : 12 Jun 2015 13:17
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC001

Quant Time: Jun 12 18:09:23 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM061215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 12 16:27:37 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	0.624	0.534	14.4	102	0.00
3	n-Nitrosodimethylamine	0.872	0.866	0.7	106	0.00
4 S	2-Fluorophenol	1.184	1.215	-2.6	113	0.00
5 S	Phenol-d6	1.488	1.554	-4.4	116	0.00
6 I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
7 S	Nitrobenzene-d5	0.383	0.405	-5.7	120	0.00
8	Nitrobenzene	0.414	0.432	-4.3	118	0.00
9	Naphthalene	1.212	1.152	5.0	108	0.00
10	Hexachlorobutadiene	0.201	0.190	5.5	109	0.00
11	2-Methylnaphthalene	0.782	0.762	2.6	111	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	110	0.00
13 S	2,4,6-Tribromophenol	0.182	0.226	-24.2	135	0.00
14 S	2-Fluorobiphenyl	1.746	1.659	5.0	111	0.00
15	Acenaphthylene	2.337	2.378	-1.8	117	-0.01
16	Acenaphthene	1.487	1.426	4.1	113	0.00
17	Fluorene	1.519	1.583	-4.2	122	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
19	4-Bromophenyl-phenylether	0.502	0.475	5.4	102	0.00
20	Hexachlorobenzene	0.455	0.487	-7.0	118	0.00
21	Pentachlorophenol	0.124	0.150	-21.0	148	0.00
22	Phenanthrene	1.640	1.409	14.1	101	0.00
23	Anthracene	1.938	2.201	-13.6	125	0.00
24	Fluoranthene	2.429	2.305	5.1	108	0.00
25 I	Chrysene-d12	1.000	1.000	0.0	93	0.00
26	Pyrene	1.716	1.844	-7.5	104	0.00
27 S	Terphenyl-d14	0.754	0.809	-7.3	103	0.00
28	Benzo(a)anthracene	1.571	1.543	1.8	95	0.00
29	Chrysene	1.503	1.450	3.5	94	0.02
30	Bis(2-ethylhexyl)phthalate	0.979	1.334	-36.3#	145	0.02
31	Indeno(1,2,3-cd)pyrene	1.181	1.182	-0.1	102	0.01
32 I	Perylene-d12	1.000	1.000	0.0	91	0.01
33	Benzo(b)fluoranthene	1.631	1.597	2.1	95	0.01
34	Benzo(k)fluoranthene	1.584	1.483	6.4	89	0.01
35 C	Benzo(a)pyrene	1.359	1.378	-1.4	98	0.01
36	Dibenzo(a,h)anthracene	1.142	1.200	-5.1	104	0.00
37	Benzo(g,h,i)perylene	1.234	1.243	-0.7	101	0.01

(#) = Out of Range

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