

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021315\
 Data File : BM000516.D
 Acq On : 13 Feb 2015 08:04
 Operator : TP/IZ
 Sample : SST0.406
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SST0.406

Quant Time: Feb 13 12:03:15 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 12 16:45:55 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	112	0.00
2 I	Naphthalene-d8	1.000	1.000	0.0	114	0.00
3	Naphthalene	0.938	0.951	-1.4	115	0.00
4 SURR	2-Methylnaphthalene-d10	0.510	0.499	2.2	111	0.00
5	2-Methylnaphthalene	0.634	0.635	-0.2	114	0.00
6 I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
7	Acenaphthylene	1.505	1.536	-2.1	116	0.00
8 C	Acenaphthene	1.155	1.186	-2.7	113	-0.02
9	Fluorene	1.396	1.419	-1.6	114	0.00
10 I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
11	Pentachlorophenol	0.051	0.049	3.9	120	-0.02
12	Phenanthrene	1.086	1.065	1.9	115	0.00
13	Anthracene	0.977	0.969	0.8	116	0.00
14 SURR	Fluoranthene-d10	0.841	0.839	0.2	114	0.00
15 C	Fluoranthene	1.211	1.228	-1.4	114	-0.01
16 I	Chrysene-d12	1.000	1.000	0.0	110	-0.01
17	Pyrene	1.439	1.464	-1.7	117	-0.01
18	Benzo(a)anthracene	1.128	1.139	-1.0	112	-0.01
19	Chrysene	1.300	1.326	-2.0	110	-0.01
20 I	Perylene-d12	1.000	1.000	0.0	105	0.00
21	Benzo(b)fluoranthene	1.352	1.335	1.3	103	-0.01
22	Benzo(k)fluoranthene	1.345	1.424	-5.9	114	-0.01
23 C	Benzo(a)pyrene	1.264	1.282	-1.4	107	-0.01
24	Indeno(1,2,3-cd)pyrene	1.410	1.350	4.3	100	-0.01
25	Dibenzo(a,h)anthracene	1.114	1.051	5.7	99	-0.01
26	Benzo(g,h,i)perylene	1.244	1.190	4.3	100	-0.01

(#) = Out of Range

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