

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE011416\
 Data File : BE091370.D
 Acq On : 15 Jan 2016 3:17
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Jan 15 15:22:36 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE011416.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 23:29:02 2016
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	121	0.00
2 S	2-Fluorophenol	1.000	1.581	-58.1#	205	-0.02
3 S	Phenol-d6	1.000	1.292	-29.2#	202	-0.01
4 I	Naphthalene-d8	5.000	5.000	0.0	132	0.00
5 S	Nitrobenzene-d5	1.000	1.057	-5.7	140	0.00
6	Naphthalene	1.000	1.020	-2.0	132	0.00
7	2-Methylnaphthalene	1.000	1.036	-3.6	178	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	142	0.00
9 S	2,4,6-Tribromophenol	1.000	1.034	-3.4	182	0.00
10 S	2-Fluorobiphenyl	1.000	1.081	-8.1	164	0.00
11	Acenaphthylene	1.000	0.995	0.5	173	0.00
12	Acenaphthene	1.000	0.961	3.9	138	0.00
13	Fluorene	1.000	0.976	2.4	157	0.00
14 I	Phenanthrene-d10	5.000	5.000	0.0	125	0.00
15	Phenanthrene	1.000	1.012	-1.2	123	-0.01
16	Anthracene	1.000	1.037	-3.7	152	0.00
17	Fluoranthene	1.000	1.215	-21.5	167	0.00
18 I	Chrysene-d12	5.000	5.000	0.0	130	0.00
19	Pyrene	1.000	1.103	-10.3	201	0.00
20 S	Terphenyl-d14	1.000	1.098	-9.8	185	0.00
21	Benzo(a)anthracene	1.000	1.207	-20.7	236	0.00
22	Chrysene	1.000	1.050	-5.0	137	0.00
23	Indeno(1,2,3-cd)pyrene	1.000	1.201	-20.1	209	0.00
24 I	Perylene-d12	5.000	5.000	0.0	141	0.00
25	Benzo(b)fluoranthene	1.000	1.166	-16.6	224	0.00
26	Benzo(k)fluoranthene	1.000	1.043	-4.3	171	0.00
27 C	Benzo(a)pyrene	1.000	1.080	-8.0	174	0.00
28	Dibenzo(a,h)anthracene	1.000	1.215	-21.5	262	0.00
29	Benzo(g,h,i)perylene	1.000	1.146	-14.6	210	0.00

(#) = Out of Range

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