

Data Path : Z:\HPCHEM1\BNA E\DATA\BE021916\
 Data File : BE091421.D
 Acq On : 19 Feb 2016 19:32
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Feb 19 22:39:42 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE020216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 00:27:53 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	92	0.00
2 S	2-Fluorophenol	1.000	1.103	-10.3	126	-0.03
3 S	Phenol-d6	1.000	1.170	-17.0	120	-0.03
4 I	Naphthalene-d8	5.000	5.000	0.0	96	-0.01
5 S	Nitrobenzene-d5	1.000	1.175	-17.5	144	-0.02
6	Naphthalene	1.000	0.936	6.4	101	-0.01
7	2-Methylnaphthalene	1.000	1.016	-1.6	113	-0.01
8 I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
9 S	2,4,6-Tribromophenol	1.000	1.054	-5.4	147	-0.01
10 S	2-Fluorobiphenyl	1.000	0.946	5.4	104	-0.01
11	Acenaphthylene	1.000	1.043	-4.3	125	0.00
12	Acenaphthene	1.000	0.914	8.6	103	0.00
13	Fluorene	1.000	0.957	4.3	104	0.00
14 I	Phenanthrene-d10	5.000	5.000	0.0	96	-0.01
15	Phenanthrene	1.000	0.929	7.1	101	0.00
16	Anthracene	1.000	1.047	-4.7	111	-0.01
17	Fluoranthene	1.000	1.054	-5.4	121	0.00
18 I	Chrysene-d12	5.000	5.000	0.0	97	0.00
19	Pyrene	1.000	1.091	-9.1	120	0.00
20 S	Terphenyl-d14	1.000	1.075	-7.5	114	0.00
21	Benzo(a)anthracene	1.000	1.120	-12.0	137	0.00
22	Chrysene	1.000	1.008	-0.8	107	0.00
23	Indeno(1,2,3-cd)pyrene	1.000	1.142	-14.2	132	0.00
24 I	Perylene-d12	5.000	5.000	0.0	106	0.00
25	Benzo(b)fluoranthene	1.000	0.990	1.0	119	0.00
26	Benzo(k)fluoranthene	1.000	1.070	-7.0	131	0.00
27 C	Benzo(a)pyrene	1.000	1.112	-11.2	140	0.00
28	Dibenzo(a,h)anthracene	1.000	1.128	-12.8	149	0.00
29	Benzo(g,h,i)perylene	1.000	1.050	-5.0	129	0.00

(#) = Out of Range

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