

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012417\
 Data File : BE092671.D
 Acq On : 24 Jan 2017 18:31
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jan 25 13:12:36 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE011617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 17 10:14:01 2017
 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	112	0.00
2	1,4-Dioxane	0.400	0.337	15.8	93	0.00
3	n-Nitrosodimethylamine	0.400	0.230	42.5#	73	0.05
4 S	2-Fluorophenol	0.400	0.285	28.8#	94	0.02
5 S	Phenol-d6	0.400	0.307	23.3	104	0.02
6	bis(2-Chloroethyl)ether	0.400	0.342	14.5	114	0.02
7 I	Naphthalene-d8	5.000	5.000	0.0	113	0.00
8 S	Nitrobenzene-d5	0.400	0.326	18.5	111	0.00
9	Naphthalene	0.400	0.389	2.8	119	0.00
10	Hexachlorobutadiene	0.400	0.392	2.0	117	0.00
11	2-Methylnaphthalene	0.400	0.370	7.5	117	0.01
12 I	Acenaphthene-d10	5.000	5.000	0.0	119	0.00
13 S	2,4,6-Tribromophenol	0.400	0.433	-8.2	135	0.03
14 S	2-Fluorobiphenyl	0.400	0.375	6.3	121	0.00
15	Acenaphthylene	0.400	0.375	6.3	126	0.00
16	Acenaphthene	0.400	0.400	0.0	126	0.00
17	Fluorene	0.400	0.378	5.5	126	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	119	0.00
19	Hexachlorobenzene	0.400	0.409	-2.2	122	0.00
20	Pentachlorophenol	0.400	0.008	98.0#	3	0.00
21	Phenanthrene	0.400	0.418	-4.5	129	0.00
22	Anthracene	0.400	0.384	4.0	129	0.00
23	Fluoranthene	0.400	0.403	-0.8	133	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	123	0.00
25	Pyrene	0.400	0.363	9.3	134	0.00
26 S	Terphenyl-d14	0.400	0.356	11.0	128	0.00
27	Benzo(a)anthracene	0.400	0.373	6.8	134	0.00
28	Chrysene	0.400	0.417	-4.2	138	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.398	0.5	92	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.358	10.5	126	0.00
31 I	Perylene-d12	5.000	5.000	0.0	121	0.00
32	Benzo(b)fluoranthene	0.400	0.387	3.3	127	0.00
33	Benzo(k)fluoranthene	0.400	0.378	5.5	133	0.00
34 C	Benzo(a)pyrene	0.400	0.370	7.5	129	0.00
35	Dibenzo(a,h)anthracene	0.400	0.366	8.5	125	0.00
36	Benzo(g,h,i)perylene	0.400	0.375	6.3	125	0.00

(#) = Out of Range

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