

Data Path : Z:\HPCHEM1\BNA E\DATA\BE122716\
 Data File : BE092617.D
 Acq On : 27 Dec 2016 10:04
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 28 01:29:27 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE121916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 19 16:16:32 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	134	0.00
2	1,4-Dioxane	0.400	0.390	2.5	146	0.00
3	n-Nitrosodimethylamine	0.400	0.482	-20.5	176	-0.01
4 S	2-Fluorophenol	0.400	0.431	-7.7	144	0.00
5 S	Phenol-d6	0.400	0.450	-12.5	164	-0.02
6	bis(2-Chloroethyl)ether	0.400	0.383	4.3	151	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	130	0.00
8 S	Nitrobenzene-d5	0.400	0.472	-18.0	153	0.00
9	Naphthalene	0.400	0.399	0.3	140	0.00
10	Hexachlorobutadiene	0.400	0.403	-0.8	140	0.00
11	2-Methylnaphthalene	0.400	0.383	4.3	139	-0.01
12 I	Acenaphthene-d10	5.000	5.000	0.0	130	0.00
13 S	2,4,6-Tribromophenol	0.400	0.435	-8.7	139	0.00
14 S	2-Fluorobiphenyl	0.400	0.389	2.8	138	0.00
15	Acenaphthylene	0.400	0.372	7.0	135	0.00
16	Acenaphthene	0.400	0.381	4.8	133	0.00
17	Fluorene	0.400	0.383	4.3	137	-0.01
18 I	Phenanthrene-d10	5.000	5.000	0.0	128	0.00
19	Hexachlorobenzene	0.400	0.419	-4.7	145	0.00
20	Pentachlorophenol	0.400	0.401	-0.3	143	0.00
21	Phenanthrene	0.400	0.370	7.5	134	0.00
22	Anthracene	0.400	0.366	8.5	138	0.00
23	Fluoranthene	0.400	0.371	7.3	135	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	129	0.00
25	Pyrene	0.400	0.365	8.8	134	0.00
26 S	Terphenyl-d14	0.400	0.370	7.5	134	0.00
27	Benzo(a)anthracene	0.400	0.388	3.0	146	0.00
28	Chrysene	0.400	0.395	1.3	138	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.371	7.3	129	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.365	8.8	137	0.00
31 I	Perylene-d12	5.000	5.000	0.0	126	0.01
32	Benzo(b)fluoranthene	0.400	0.374	6.5	135	0.00
33	Benzo(k)fluoranthene	0.400	0.397	0.8	144	0.00
34 C	Benzo(a)pyrene	0.400	0.376	6.0	141	0.00
35	Dibenzo(a,h)anthracene	0.400	0.361	9.8	136	0.02
36	Benzo(g,h,i)perylene	0.400	0.378	5.5	135	0.00

(#) = Out of Range

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