

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081116\
 Data File : BE092267.D
 Acq On : 11 Aug 2016 15:15
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 11 17:11:48 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 10 11:12:32 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	146	0.00
2	1,4-Dioxane	0.400	0.378	5.5	136	0.00
3	n-Nitrosodimethylamine	0.400	0.475	-18.7	161	-0.01
4 S	2-Fluorophenol	0.400	0.398	0.5	160	0.00
5 S	Phenol-d6	0.400	0.381	4.8	178	-0.02
6	bis(2-Chloroethyl)ether	0.400	0.431	-7.7	175	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	145	0.00
8 S	Nitrobenzene-d5	0.400	0.376	6.0	161	-0.02
9	Naphthalene	0.400	0.398	0.5	145	-0.02
10	Hexachlorobutadiene	0.400	0.403	-0.8	147	-0.02
11	2-Methylnaphthalene	0.400	0.401	-0.3	148	-0.01
12 I	Acenaphthene-d10	5.000	5.000	0.0	143	0.00
13 S	2,4,6-Tribromophenol	0.400	0.497	-24.2	158	0.00
14 S	2-Fluorobiphenyl	0.400	0.394	1.5	145	-0.01
15	Acenaphthylene	0.400	0.389	2.8	145	0.00
16	Acenaphthene	0.400	0.397	0.8	144	0.00
17	Fluorene	0.400	0.394	1.5	144	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	145	0.00
19	Hexachlorobenzene	0.400	0.363	9.3	131	0.00
20	Pentachlorophenol	0.400	0.441	-10.2	186	0.00
21	Phenanthrene	0.400	0.374	6.5	140	0.00
22	Anthracene	0.400	0.376	6.0	147	0.00
23	Fluoranthene	0.400	0.374	6.5	143	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	147	-0.02
25	Pyrene	0.400	0.408	-2.0	157	0.00
26 S	Terphenyl-d14	0.400	0.406	-1.5	161	0.00
27	Benzo(a)anthracene	0.400	0.433	-8.2	171	-0.02
28	Chrysene	0.400	0.393	1.8	142	0.00
29	Bis(2-ethylhexyl)phthalate	0.400	0.490	-22.5	157	0.00
30	Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	164	-0.02
31 I	Perylene-d12	5.000	5.000	0.0	148	0.00
32	Benzo(b)fluoranthene	0.400	0.398	0.5	160	0.00
33	Benzo(k)fluoranthene	0.400	0.407	-1.7	150	0.00
34 C	Benzo(a)pyrene	0.400	0.409	-2.2	166	0.00
35	Dibenzo(a,h)anthracene	0.400	0.404	-1.0	167	0.00
36	Benzo(g,h,i)perylene	0.400	0.392	2.0	153	-0.02

(#) = Out of Range

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