

Data Path : Z:\HPCHEM1\BNA E\Data\BE041116\
 Data File : BE091633.D
 Acq On : 11 Apr 2016 18:02
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 12 15:35:55 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 16:48:14 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	120	0.00
2	1,4-Dioxane	0.400	0.370	7.5	114	0.00
3	n-Nitrosodimethylamine	0.400	0.359	10.3	116	0.00
4 S	2-Fluorophenol	0.400	0.341	14.8	108	0.00
5 S	Phenol-d6	0.400	0.346	13.5	114	0.00
6	bis(2-Chloroethyl)ether	0.400	0.380	5.0	120	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	122	-0.02
8 S	Nitrobenzene-d5	0.400	0.329	17.8	114	0.00
9	Nitrobenzene	0.400	0.432	-8.0	113	0.00
10	Naphthalene	0.400	0.396	1.0	123	0.00
11	Hexachlorobutadiene	0.400	0.381	4.8	118	0.00
12	2-Methylnaphthalene	0.400	0.388	3.0	123	-0.01
13 I	Acenaphthene-d10	5.000	5.000	0.0	123	0.00
14 S	2,4,6-Tribromophenol	0.400	0.452	-13.0	111	0.00
15 S	2-Fluorobiphenyl	0.400	0.391	2.3	123	0.00
16	Acenaphthylene	0.400	0.359	10.3	137	0.00
17	Acenaphthene	0.400	0.409	-2.2	131	-0.01
18	Fluorene	0.400	0.382	4.5	122	-0.01
19 I	Phenanthrene-d10	5.000	5.000	0.0	122	0.00
20	4-Bromophenyl-phenylether	0.400	0.374	6.5	118	0.00
21	Hexachlorobenzene	0.400	0.397	0.8	122	0.00
22	Pentachlorophenol	0.400	0.421	-5.2	126	0.00
23	Phenanthrene	0.400	0.408	-2.0	125	0.00
24	Anthracene	0.400	0.379	5.3	123	0.00
25	Fluoranthene	0.400	0.367	8.3	121	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	114	0.00
27	Benzydine	0.400	0.518	-29.5#	124	-0.02
28	Pvrene	0.400	0.384	4.0	119	-0.02
29 S	Terphenyl-d14	0.400	0.432	-8.0	126	0.00
30	Benzo(a)anthracene	0.400	0.392	2.0	116	0.00
31	3,3'-Dichlorobenzidine	0.400	0.340	15.0	87	0.00
32	Chrysene	0.400	0.454	-13.5	132	-0.02
33	Bis(2-ethylhexyl)phthalate	0.400	0.345	13.8	63	0.00
34	Indeno(1,2,3-cd)pyrene	0.400	0.348	13.0	103	-0.01
35 I	Perylene-d12	5.000	5.000	0.0	111	-0.01
36	Benzo(b)fluoranthene	0.400	0.363	9.3	106	-0.01
37	Benzo(k)fluoranthene	0.400	0.404	-1.0	118	-0.01
38 C	Benzo(a)pyrene	0.400	0.344	14.0	102	0.00
39	Dibenzo(a,h)anthracene	0.400	0.353	11.8	103	-0.01
40	Benzo(a,h,i)perylene	0.400	0.360	10.0	104	-0.01

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Compound	Amount	Calc.	%Dev Area	% Dev(min)
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(#) = Out of Range

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