

## Getting data in and getting results out

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# DATA\_SECTION

- Is where data is read in
- If needed it can be processed a bit before entering the likelihood
- This section is only evaluated ones
- AD Model Builder will not keep track of the derivatives for the quantities declared here
- We have already seen a few examples

```
# number of observations
10
# observed Y values
1.4 4.7 5.1 8.3 9.0 14.5 14.0 13.4 19.2 18
# observed x values
-1 0 1 2 3 4 5 6 7 8

DATA_SECTION
  init_int N
  init_vector Y(1,N)
  init_vector x(1,N)

PARAMETER_SECTION
  init_number a
  init_number b
  init_number logSigma
  sdreport_number sigmasq
  objective_function_value nll

PROCEDURE_SECTION
  sigmasq=exp(2*logSigma);
  nll=0.5*(N*log(2*M_PI*sigmasq)+sum(square(Y-(a+b*x)))/sigmasq);
```

# Basic rules

- Lines in the data file starting with a **#** are comments and ignored by AD Model Builder
- All types in the DATA\_SECTION starting with **init\_** are initialized from the data file
- E.g:

**init\_int**

**init\_number**

**init\_ivector**

**init\_vector**

**init\_imatrix**

**init\_matrix**

**init\_3darray ... init\_7darray**

**init\_adstring**

- All types without **init\_** can be used for further calculations, but are not initialized from the data file (unless it is done ‘manually’)

# Dimensions set by data

- Notice that the length of the vectors **N** are also read in
- This is good when we want to make general programs, as no recompilation is needed to run the same program for a data set of different length

```
# number of observations
10
# observed Y values
1.4 4.7 5.1 8.3 9.0 14.5 14.0 13.4 19.2 18
# observed x values
-1 0 1 2 3 4 5 6 7 8
```

```
DATA_SECTION
init_int N
init_vector Y(1,N)
init_vector x(1,N)
```

# Data matrix and local one line calculations

- A different way to input data for the same example

```
DATA_SECTION
  init_int N
  init_matrix xy(1,N,1,2)

  vector x(1,N)
  vector Y(1,N)

  !! x = column(xy,1);
  !! Y = column(xy,2);
```

	# number of observations
	10
	# observed x and Y values
-1	1.4
0	4.7
1	5.1
2	8.3
3	9.0
4	14.5
5	14.0
6	13.4
7	19.2
8	18

- First a matrix is read from the data file
- Then the first column is saved as a vector 'x' and the second column as a vector 'Y'
- The rest of the program is unchanged
- A common use for this feature is transformations like:

```
DATA_SECTION
  init_int N
  init_vector obs(1,N)

  vector logObs(1,N)
  !! logObs = log(obs);
```

# LOC\_CALCS and random numbers

- For longer calculations **LOC\_CALCS . . . END\_CALCS** is more convenient than  
**!!**

```
DATA_SECTION
  vector X(1,1000);
LOC_CALCS
  random_number_generator rng(123456);
  X.fill_randn(rng);
  X*=5.0;
  X+=2.0;
END_CALCS
PARAMETER_SECTION
  init_number logSigma;
  init_number mu;
  sdreport_number sigma;
  objective_function_value nll;
PROCEDURE_SECTION
  sigma=exp(logSigma);
  int N=X.indexmax()-X.indexmin()+1;
  dvariable ss=square(sigma);
  nll=0.5*(N*log(2*M_PI*ss)+sum(square(X-mu))/ss);
```

- This program uses simulated data only, so the data file is not needed, except ...

# More on generating random numbers

```
DATA_SECTION
LOC_CALCS
  random_number_generator rng(123456);
  dvector sample(1,5);

  sample.fill_randu(rng);
  cout<<"Uniform(0,1): "<<sample<<endl;

  sample.fill_randn(rng);
  cout<<"Normal(0,1): "<<sample<<endl;

  sample.fill_randpoisson(1.5,rng);
  cout<<"pois(1.5): "<<sample<<endl;

  sample.fill_randnegbinomial(1.5,2.0,rng);
  cout<<"neg.bin(1.5,2): "<<sample<<endl;

  sample.fill_randcau(rng);
  cout<<"Cauchy: "<<sample<<endl;

  sample.fill_randbi(0.8,rng);
  cout<<"binomial(n=1,p=0.8): "<<sample<<endl;

  dvector p("{.01,.495,.495}");
  sample.fill_multinomial(rng,p);
  cout<<"multinomial(n=1,p=(.01,.495,.495)): "
    <<sample<<endl;

  ad_exit(0);
END_CALCS
PARAMETER_SECTION
  objective_function_value nll;

PROCEDURE_SECTION
```

```
Uniform(0,1):  0.779837 0.229835 0.0126429
               0.714228 0.654815

Normal(0,1):  -0.325127 1.03682 0.567672
               -0.670345 2.89024

pois(1.5):    2 2 0 1 2

neg.bin(1.5,2):  0 3 1 0 4

Cauchy:  -0.188267 -1.30511 -9.11156
          29.8652  2.83259

binomial(n=1,p=0.8):  1 1 0 0 1

multinomial(n=1,p=(.01,.495,.495)):  3 2 3 2 3
```

# Changing input file

- The default is to read from a file named **<modelname>.dat**, where <modelname> is the name of the \*.tpl file
- If for some reason we feel like reading from another file we can run the program with the command line option

```
an@ch-pcb-an:~/$. /model -ind newdatafile.dat
```

- Or within the DATA\_SECTION use a command like

```
DATA_SECTION
  init_int nrowA
  init_int ncolA
  init_matrix A(1,nrowA,1,ncolA)

  !! ad_comm::change_datafile_name("newfile.dat");

  init_int nrowB
  init_int ncolB
  init_matrix B(1,nrowB,1,ncolB)
```

- Then A is read from the default file and B are read from 'newfile.dat'



# Ragged arrays

- A special feature is to use integer vectors (ivector) as dimensions to get ‘ragged arrays’

DATA_SECTION	
init_int N	2
init_ivector startYear(1,N)	1991 1995
init_ivector endYear(1,N)	2000 1999
init_matrix A(1,N,startYear,endYear)	23 5 54 12 45 8 23 45 76 32
	45 34 32 54 34

- Notice that the two rows in ‘A’ does not have the same column index, and that is OK

# Checking what is read in

- John once taught me a neat way to check the input, consider this:

```
GLOBALS_SECTION
#include <fstream.h>
ofstream clogf("program.log");
#define TRACE(obj) clogf<<"line "<<__LINE__<<", file "<<__FILE__<<", "<<#obj" =\n " \
<<obj<<endl<<endl;

DATA_SECTION
init_int N
!! TRACE(N)
init_ivector startYear(1,N)
!! TRACE(startYear)
init_ivector endYear(1,N)
!! TRACE(endYear)
init_matrix A(1,N,startYear,endYear)
!! TRACE(A)
!! ad_exit(0);
PARAMETER_SECTION
objective_function_value nll;
PROCEDURE_SECTION
```

- Then the following is placed in the ‘program.log’, and we can easily check it

```
line 15, file ra.cpp, N =
2

line 17, file ra.cpp, startYear =
1991 1995

line 19, file ra.cpp, endYear =
2000 1999

line 21, file ra.cpp, A =
23 5 54 12 45 8 23 45 76 32
45 34 32 54 34
```

# Reading standard results into R

- The following R-function is useful for reading the standard ADMB output files

```
read.fit<-function(file){
  #
  # Function to read a basic AD Model Builder fit.
  #
  # Use for instance by:
  #
  #   simple.fit <- read.fit('c:/admb/examples/simple')
  #
  # Then the object 'simple.fit' is a list containing sub-objects
  # 'names', 'est', 'std', 'cor', and 'cov' for all model
  # parameters and sdreport quantities.
  #
  ret<-list()
  parfile<-as.numeric(scan(paste(file, '.par', sep=''),
                           what='', n=16, quiet=TRUE)[c(6,11,16)])
  ret$npar<-as.integer(parfile[1])
  ret$nlogl<-parfile[2]
  ret$maxgrad<-parfile[3]
  file<-paste(file, '.cor', sep='')
  lin<-readLines(file)
  ret$npar<-length(lin)-2
  ret$logDetHess<-as.numeric(strsplit(lin[1], '=')[[1]][2])
  sublin<-lapply(strsplit(lin[1:ret$npar+2], ' '),function(x)x[x!=''])
  ret$names<-unlist(lapply(sublin,function(x)x[2]))
  ret$est<-as.numeric(unlist(lapply(sublin,function(x)x[3])))
  ret$std<-as.numeric(unlist(lapply(sublin,function(x)x[4])))
  ret$cor<-matrix(NA, ret$npar, ret$npar)
  corvec<-unlist(sapply(1:length(sublin), function(i)sublin[[i]][5:(4+i)]))
  ret$cor[upper.tri(ret$cor, diag=TRUE)]<-as.numeric(corvec)
  ret$cor[lower.tri(ret$cor)] <- t(ret$cor)[lower.tri(ret$cor)]
  ret$cov<-ret$cor*(ret$std%o%ret$std)
  return(ret)
}
```

```

> source("tools.R")
> fit <- read.fit("simplelm")
> fit

$nopar
[1] 3

$nlogl
[1] 17.6406

$maxgrad
[1] 1.0698e-06

$npar
[1] 4

$logDetHess
[1] 8.33061

$names
[1] "a"          "b"          "logSigma" "sigmasq"

$est
[1] 4.07820 1.90910 0.34513 1.99420

$std
[1] 0.70394 0.15547 0.22361 0.89184

$cor
      [,1] [,2] [,3] [,4]
[1,] 1.000 -0.773  0    0
[2,] -0.773 1.000  0    0
[3,] 0.000 0.000  1    1
[4,] 0.000 0.000  1    1

$cov
      [,1] [,2] [,3] [,4]
[1,] 0.49553152 -0.08459832 0.00000000 0.00000000
[2,] -0.08459832 0.02417092 0.00000000 0.00000000
[3,] 0.00000000 0.00000000 0.05000143 0.1994243
[4,] 0.00000000 0.00000000 0.19942434 0.7953786

```

# The REPORT\_SECTION

- The REPORT\_SECTION is for user defined output.
- Any calculated quantity can be written, and formatted as desired
- For instance:

```
REPORT_SECTION  
report<<"Here comes the matrix X"<<endl;  
report<<X<<endl;  
report<<"After that the vector y"<<endl;  
report<<y<<endl;
```

- The output is written to the file `<modelname>.rep`

# Exercises

## Exercise 1: Inputting and using prior information in a Beverton-Holt model

- The Beverton-Holt model can be written (slightly re-parametrized) as:

$$\log R = \log(a) + \log(ssb) - \log(1 + \exp(\log(b))ssb)$$

- We want to estimate the model parameters  $\log(a)$  and  $\log(b)$  and have two sources.
- A data set of SSB and  $\log(R)$

<http://www.nielsensweb.org/TCADSA/dat/BHex/bh.dat>

- Which can be modeled with a program like

<http://www.nielsensweb.org/TCADSA/dat/BHex/bh.tpl>

- The second source is the result of for a similar species in a similar area.

1	loga	1.8085e+00	1.2725e-01	1.0000	
2	logb	-1.2183e+01	3.2431e-01	0.9278	1.0000

- Modify the program to use this prior information on  $\log(a)$  and  $\log(b)$

## Solution:

```
GLOBALS_SECTION
#include <fstream.h>
ofstream clogf("program.log");
#define TRACE(obj) clogf<<"line "<<__LINE__<<", file "<<__FILE__<<", "<<#obj" =\n " \
    <<obj<<endl<<endl;

DATA_SECTION
init_int nR
init_int nC
init_matrix obs(1,nR,1,nC)
vector ssb(1,nR)
!! ssb=column(obs,1);
vector logR(1,nR)
!! logR=column(obs,2);

!! ad_comm::change_datafile_name("prior.cor");

init_int idxA
init_adstring nameA
init_number estA
init_number sdA
init_number corrAA

init_int idxB
init_adstring nameB
init_number estB
init_number sdB
init_number corrBA
init_number corrBB

vector meanAB(1,2)
matrix covAB(1,2,1,2)
LOC_CALCS
meanAB(1)=estA;
meanAB(2)=estB;
covAB(1,1)=square(sdA);
covAB(1,2)=sdA*sdB*corrBA;
covAB(2,1)=covAB(1,2);
covAB(2,2)=square(sdB);
TRACE(meanAB);
TRACE(covAB);
END_CALCS
PARAMETER_SECTION
init_number loga;
```

```

init_number logb;
init_number logSigma;
number sigmaSq;
vector pred(1,nR);
vector vecAB(1,2);

objective_function_value nll;

PROCEDURE_SECTION
sigmaSq=exp(2.0*logSigma);
pred=loga+log(ssb)-log(1+exp(logb)*ssb);
nll=0.5*(nR*log(2*M_PI*sigmaSq)+sum(square(logR-pred))/sigmaSq);

vecAB(1)=loga;
vecAB(2)=logb;
dvar_vector diff=vecAB-meanAB;
nll+=0.5*(log(2.0*M_PI)*2.0+log(det(covAB))+diff*inv(covAB)*diff);

```