

AD Model Builder introduction course

Data input and reporting results

AD Model Builder foundation

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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 \mathbf{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_rndn(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_ranxn(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  
    init_ivector startYear(1,N)  
    init_ivector endYear(1,N)  
    init_matrix A(1,N,startYear,endYear)
```

2	1991 1995
	2000 1999
	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$nopar<-as.integer(parfile[1])  
  ret$nlogl<-parfile[2]  
  ret$maxgrad<-parfile[3]  
  file<-paste(file,'.cor', sep='')  
  lin<-readLines(file)  
  ret$npars<-length(lin)-2  
  ret$logDetHess<-as.numeric(strsplit(lin[1], '=')[[1]][2])  
  sublin<-lapply(strsplit(lin[1:ret$npars+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$npars, ret$npars)  
  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

$npars
[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/ADMB2/BHex/bh.dat>

- Which can be modeled program like

<http://www.nielsensweb.org/ADMB2/BHex/bh.tpl>

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

```
The logarithm of the determinant of the hessian = 12.869
index    name        value      std dev     1       2       3
  1    loga      1.8085e+00 1.2725e-01  1.0000
  2    logb      -1.2183e+01 3.2431e-01  0.9278  1.0000
  3  logSigma   -1.1332e+00 1.0426e-01  0.0000  0.0000  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");
!! // Dirty trick to wind fast forward over the first two lines
!! // Alternatively the two lines could have been commented out with #
!! for(int i=1; i<=18; ++i){
init_adstring tmp;
!!
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);

```