

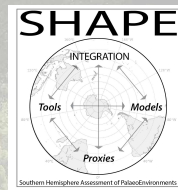
Modelling using CSIRO Mk3L

Part 2: Design your own experiment

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1. Working with model output

Output files

- When the model runs, it generates output. This is what you want!
- The model generates two types of output:

output files save the state of the model *during* a simulation

restart files save the state of the model at the *end* of a simulation

- The output files contain the simulated climate.
- In common with almost all climate models, CSIRO Mk3L saves its output in a format called netCDF.
- netCDF is a self-describing, machine-independent data format. For further information see:
 - <http://www.unidata.ucar.edu/software/netcdf/>

Ferret

- A free data visualisation and analysis package.
- Specifically designed for visualising climatic data.
- Makes it a breeze to visualise, analyse and manipulate the contents of netCDF files.
- Very powerful and easy-to-use averaging, interpolation and re-gridding capabilities.
- Your new best friend!
- For further information see:
 - <http://ferret.pmel.noaa.gov/Ferret/>

Basic Ferret commands

<code>use <file></code>	Load the netCDF file <file>
<code>show data</code>	List the data which is available
<code>list <variable></code>	List the values of <variable>
<code>plot <variable></code>	Produce a line plot of <variable>
<code>shade <variable></code>	Produce a shade plot of <variable>
<code>fill <variable></code>	Produce a filled plot of <variable>
<code>contour <variable></code>	Produce a contour plot of <variable>
<code>exit or q</code>	Exit

Basic Ferret transformations

- If the variable `tsc` contains surface air temperature as a function of longitude and latitude, then you can slice and dice the data using these expressions:

<code>tsc[i=10,j=8]</code>	Temperature at gridpoint (10, 8)
<code>tsc[x=140e,y=35s]</code>	Temperature at 140°E, 35°S
<code>tsc[x=90e:180e,y=45s:0]</code>	Temperature within 90–180°E, 45–0°S
<code>tsc[i=@ave]</code>	Zonal-mean temperature
<code>tsc[i=@ave,j=@ave]</code>	Global-mean temperature
<code>tsc[i=@max,j=@max]</code>	Global-maximum temperature
<code>tsc[i=@min,j=@min]</code>	Global-minimum temperature

Exercise 1: Ferret and model output

- First, get the course material:

```
cd ~/<NAME>
tar zxvf /srv/scratch/z3210932/material.tar.gz
```

- These commands create a new directory, `material`, which contains some material for this course. This includes some typical output from a CSIRO Mk3L simulation.
- Change to this directory by entering the command:

```
cd material
```


Exercise 1: Ferret and model output

- Now, load and run Ferret:

```
module load ferret
ferret
```

- Within Ferret, load the sample atmosphere model output:

```
yes? use stsc_spi62.nc
```

- This file contains data for surface air temperature.

Exercise 1: Ferret and model output

- Try commands such as:

```
show data
fill tsc[k=1,l=1]
fill tsc[k=@ave,l=@ave]
fill tsc[i=@ave,k=@ave]
fill tsc[k=@max,l=@max]
plot tsc[i=@ave,j=@ave,k=@ave]
plot tsc[i=@ave,k=@ave,l=@ave]
plot tsc[x=140e,y=35s,l=@ave]
list tsc[i=@ave,j=@ave,k=@ave,l=@ave]
show transform
```

Exercise 1: Ferret and model output

- A sample ocean model output file is also provided. Within Ferret, load this file:

```
yes? use com.spi62.00001.nc
```

- Try commands such as:

```
shade/lev=1d temp[k=1,l=1]  
fill/lev=1d temp[i=@ave,l=@ave]  
fill/lev=2dc motg[l=@ave]  
plot mota[y=30n:60n@max,k=@max]
```

2. Configuring CSIRO Mk3L

Running and configuring CSIRO Mk3L

- The three basic steps involved in running the model are:
 - Create a run directory
 - Copy everything that you need to this directory
 - Run the model
- The “everything” in this second step consists of:
 - The model itself (the “executable”)
 - All the input files needed to run the model
- To configure the model for a particular experiment, we need to modify one or more of these input files.

Input files

- The model requires three types of input files:

control file	configures the model for a particular simulation
restart files	initialise the model at the <i>start</i> of a simulation
auxiliary files	provide the boundary conditions <i>during</i> a simulation

- The model may be configured for a particular scenario by modifying one or more of these files.
- Auxiliary files provide the boundary conditions that the model cannot simulate itself e.g. topography.
- See Chapters 4 and 5 of the Users Guide for further information.

The control file

- To run the model, you use a command such as:

```
./model < input > output
```

- The file `input` is the *control file*.
- This file contains a number of `namelist` groups.
- The parameters contained within these groups specify:
 - the duration of a simulation
 - the physical configuration of the model
 - which model variables are to be saved

namelist groups

- A namelist group looks like this:

```
&control  
  lcouple=T  
  locean=F  
  mstep=20  
  nsstop=0  
  ndstop=1  
  lastmonth=0  
  months=0  
  nrad=6  
&end
```


nano

- nano is a simple Linux text editor.
- To edit a file, enter the command:

```
nano <file>
```

- Some basic nano commands are:

```
Ctrl-G  Get Help  
Ctrl-O  Write (Save)  
Ctrl-X  Exit
```

Exercise 2: Editing control files

- Change to this directory, which contains some sample control files:

```
cd ~/<NAME>/version-1.2/core/control
```

- Create a copy of one of the control files, using a command such as:

```
cp input_cp1_1day input_copy
```

- Use `nano` to examine and edit this file.

Basic namelist options

`nsstop`, `ndstop`, `lastmonth`, `months`

These determine the duration of the simulation:

<code>nsstop</code>	Stop after <code>nsstop</code> timesteps
<code>ndstop</code>	Stop after <code>ndstop</code> days
<code>lastmonth</code>	Stop at the end of calendar month <code>lastmonth</code> (1=January, 2=February, ..., 12=December)
<code>months</code>	Stop after <code>months</code> months

The first of these to have a non-zero value is the one that takes effect.

Basic namelist options

bpyear, csolar

- bpyear specifies the epoch, in years before present (where the “present” is the year 1950 CE)
- csolar specifies the solar constant, in Wm^{-2}

runtype

- runtype specifies the name of the experiment

Exercise 3: Basic namelist options

- Look at the control files in the following directories:

```
~/<NAME>/version-1.2/core/control
```

```
~/<NAME>/material/exp01
```

```
~/<NAME>/material/exp02
```

```
~/<NAME>/material/exp04
```

- Find the following parameters, and see how the values differ:

```
nsstop, ndstop, lastmonth, months
```

```
bpyear, csolar
```

```
runtype
```

3. Using auxiliary files

The control file versus auxiliary files

- Using the control file, we can specify:
 - the duration of a simulation
 - the physical configuration of the model
 - which model variables are to be saved
- However, to configure other aspects of the model we need to modify the auxiliary files. Examples include:
 - Topography
 - Bathymetry
 - Albedo
 - Vegetation and soil types
 - CO₂ transmission coefficients
 - Ozone mixing ratios

Changing the atmospheric CO₂ concentration

- The CO₂ transmission coefficients are read from an auxiliary file.
- These files are generated by the utility `radint`.
- To compile and initialise this utility, change to the directory:

```
cd ~/<NAME>/version-1.2/pre/co2
```

- Then enter the commands:

```
make  
./pset -n 18
```

Changing the atmospheric CO₂ concentration

- To generate the auxiliary file for an atmospheric CO₂ concentration of <concentration> ppm, enter the command:

```
./radint -c <concentration>
```

- For example, for a CO₂ concentration of 280 ppm:

```
./radint -c 280
```

- This generates a file called co2_data.

Exercise 4: Changing the atmospheric CO₂ concentration

- Compile and initialise `radint` by entering the commands:

```
cd ~/<NAME>/version-1.2/pre/co2
make
./pset -n 18
```

- Now generate auxiliary files for CO₂ concentrations of 280, 560 and 1120 ppm e.g.

```
./radint -c 280
```

- Remember to rename the auxiliary file each time e.g.

```
mv co2_data co2_data.280ppm
```

Applying freshwater hosing

- To apply freshwater hosing, use these `namelist` parameters:

```
hosing_flag  If T, apply freshwater hosing
hosing_rate  The freshwater hosing rate (Sv)
```

- You must also supply the auxiliary file `hosemask`.
- A sample auxiliary file is provided with the model:

```
~/<NAME>/version-1.2/core/data/atmosphere/hosing/hosemask
```

Exercise 5: Design your own freshwater hosing mask

- Change to the directory containing the sample freshwater hosing mask and create your own copy e.g.

```
cd ~/<NAME>/version-1.2/core/data/atmosphere/hosing
cp hosemask hosemask_new
```

- Now use nano to edit this file:

```
nano hosemask_new
```

- The number 7 indicates land: don't change these values!
- Put 1 where you want the water to go, and 0 everywhere else.

4. Design your own experiment

Exercise 6a: Pre-configured experiments

- Now it's time to design and run your first experiment!
- If you're new to climate modelling, then I recommend that you choose one of the following experiments:

exp01 Control simulation

exp02 Mid-Holocene (6,000 years BP)

exp03 Last Glacial Maximum (21,000 years BP)

exp04 Snowball Earth

exp05 $2\times\text{CO}_2$

exp06 Water hosing

- Each of these experiments has already been set up for you.

Exercise 6a: Pre-configured experiments

- For your experiment, change to the appropriate directory e.g.

```
cd ~/<NAME>/material/exp01
```

- Now start your experiment e.g.

```
qsub qsub_exp01
```

- Look at the script which carries out each experiment.
- How does it differ from the control simulation (exp01)?
- What would you change if you wanted to run your experiment for 50 years, rather than 10 years?

Exercise 6b: Design your own experiment

- If you're a climate modelling guru, then design and run your own experiment instead.
- Ideas:
 - Change the epoch: simulate the past or future
 - Change the solar constant: $\pm 5\%$, $\pm 10\%$, more?
 - Change the CO₂ concentration: $\times \frac{1}{2}$, $\times 2$, $\times 4$?
 - Freshwater hosing: melting of the Greenland or Antarctic ice sheets
- Tips:
 - Use one of the pre-configured experiments as a basis.
 - Unless you're feeling extremely confident, get me to check your experiments before you run them.
 - Once you've started your experiment, use `qstat` to monitor progress.