

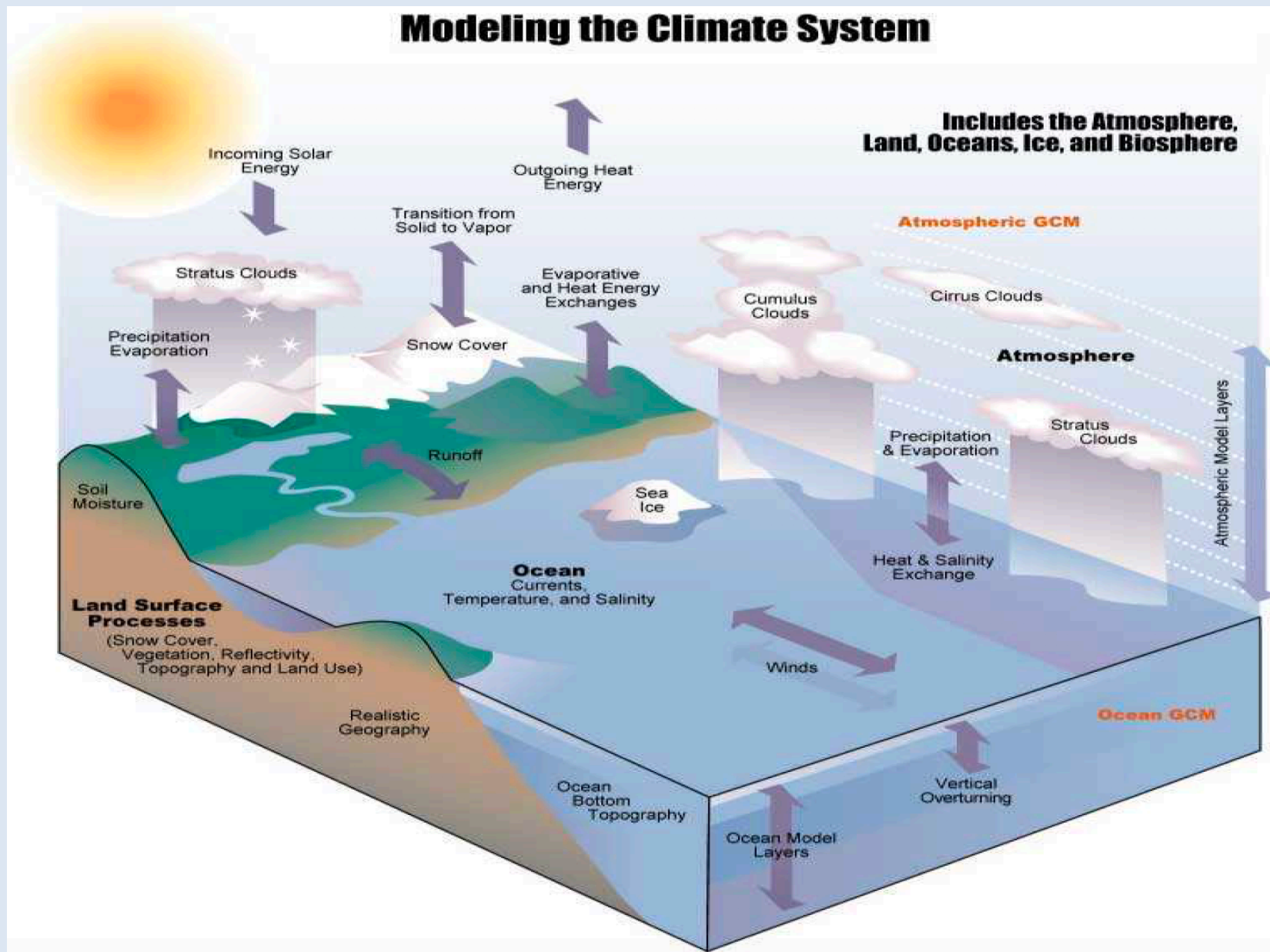
# Using the CSIRO Mk3L climate system model

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# Modeling the Climate System

**Includes the Atmosphere,  
Land, Oceans, Ice, and Biosphere**



# Choosing the right model for you

- A model is a tool - the type that you use depends upon the question that you want to answer.
- Which components, processes or quantities of the climate system do you need to model?
- Do you need a regional or a global model?
- How much spatial resolution do you need?
- How long do you need to run the model for? (for example, it isn't feasible to run a high-resolution global model for 10,000 years)
- No model is a perfect representation of the real world.

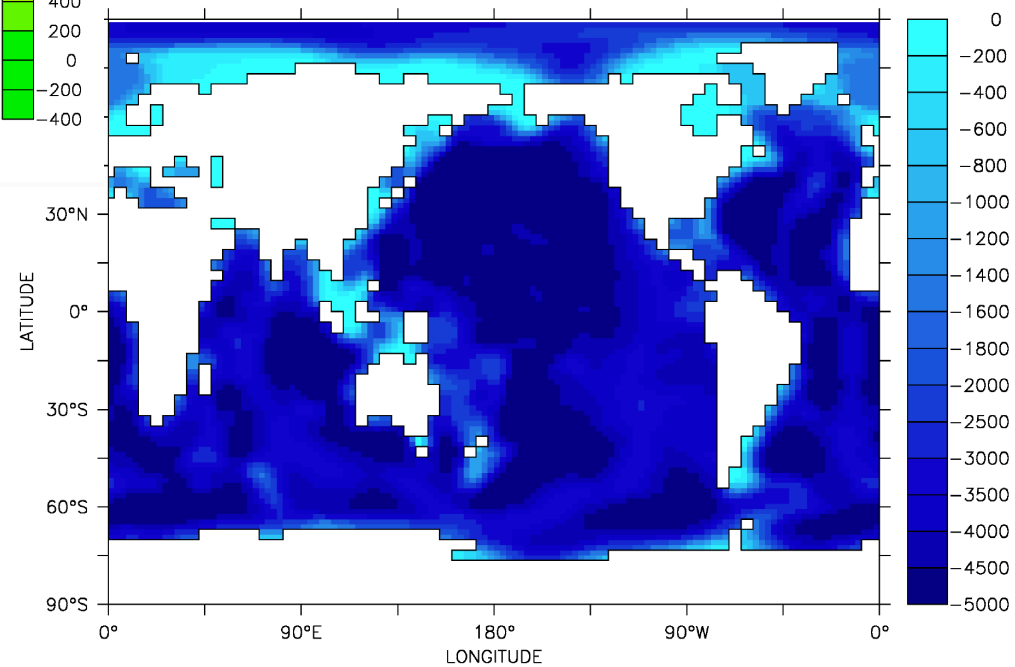
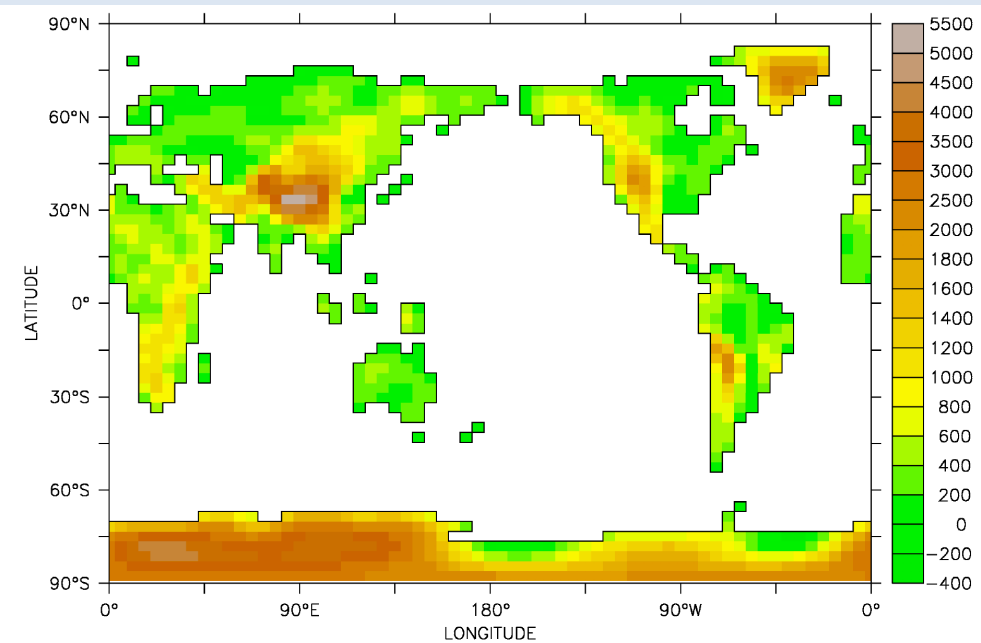
# The CSIRO Mk3L climate system model

- Low-resolution version of the CSIRO climate system model (e.g. IPCC 1st, 2nd, 3rd, 4th and 5th Assessment Reports).
- Coupled atmosphere-land-sea ice-ocean general circulation model.
- Designed to enable millennial-scale simulations of climate variability and change
- Can simulate 1000 years in around a month.
- Community model.

# The CSIRO Mk3L climate system model

- Atmosphere:
  - Three-dimensional general circulation model
  - Horizontal resolution of 5.6 3.2 with 18 vertical levels
- Ocean:
  - Three-dimensional general circulation model
  - Horizontal resolution of 2.8 1.6 with 21 vertical levels
- Sea ice:
  - Dynamic-thermodynamic sea ice model
  - Three layers (two ice, one snow)
- Land surface:
  - Soil-canopy scheme (13 land surface/vegetation types, 9 soil types)
  - Six soil layers, three snow layers

# The topography and bathymetry within CSIRO Mk3L



# Using Raijin

- Launch Terminal
- Connect through secure shell into Raijin while enabling for pop up figures
  - In the terminal, enter `userid@raijin.nci.org.au`
  - Type yes to allow login
  - Type your password
- Familiarise yourself with the basic Linux commands (see the next slide).

# Basic Linux commands

<code>ls</code>	list the contents of a directory
<code>ls -l</code>	create a long listing
<code>mkdir &lt;directory&gt;</code>	create the directory <directory>
<code>cd &lt;directory&gt;</code>	change to the directory <directory>
<code>cp &lt;file1&gt; &lt;file2&gt;</code>	copy the file <file1> to <file2>
<code>mv &lt;file1&gt; &lt;file2&gt;</code>	move the file <file1> to <file2>
<code>rm &lt;file&gt;</code>	delete the file <file>
<code>rmdir &lt;directory&gt;</code>	delete the directory <directory>
<code>man &lt;command&gt;</code>	display the manual page for <command>

- For some more Linux commands see:  
[www.dummies.com/how-to/content/linux-for-dummies-cheat-sheet.html](http://www.dummies.com/how-to/content/linux-for-dummies-cheat-sheet.html)



# Getting CSIRO Mk3L

- Mk3L is available via "subversion" but we're not going to use this today.
- To save time, there's a copy of the model distribution set as a module on Raijin.
- Install the version 1.2 of CSIRO Mk3L module by entering these commands:

```
module use /g/data/hh5/public/modules
```

```
module load mk3l/1.2
```

# Running CSIRO Mk3L

- The basic command which runs Mk3L is simply:

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

- `model` is the executable. This is the “model”.
  - `input` is the control file. This contains the instructions which tell
  - the model what to do.
  - `output` contains diagnostic information generated by the model.
- 
- The above command executes the model and feeds it the information contained within the control file and redirects the diagnostic information to an output file.

# Running CSIRO Mk3L

- Production jobs can take weeks or months to complete, we therefore need to use a queueing system.
- Raijin uses the Portable Batch System (PBS).
- Run the model by entering this command:

```
qsub qsub_test_cpl
```

- This runs the model for one day using the `qsub_test_cpl` script which launches the model for a 1 day run.
- The `qsub` command submits a job to the queueing system.
- The file `qsub_test_cpl` tells the queueing system what to do.
- Use the command `nqstat` to check the progress of your job:

```
nqstat
```

# Running CSIRO Mk3L

- The file `qsub_test_cp1` is called a script. The instructions contained within this file describe how to run the model.

- Using the `vi` command, examine the contents:

```
vi qsub_test_cp1
```

- Hint: type `ESC`, then `q!` to exit `vi`.
- The lines beginning with `#` are comments.
- The lines beginning with `#PBS -l` tell the queueing system which resources are required to run the job.

# Requesting resources

- When using a queueing system, you need to request sufficient resources to run your job.
- The script that you just ran uses three different options to do this:

<code>nodes</code>	The number of nodes to run on
<code>vmem</code>	The total amount of memory required
<code>walltime</code>	The expected run time

- It's important to request sufficient resources, but not too much.
- For further information see:  
<https://nci.org.au/user-support/training/training-exercises/parallel-programming/>

# 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.
- See Chapter 6 of the Users Guide for further information.
- In common with almost all climate models, CSIRO Mk3L saves its output in a format called netCDF.

# netCDF

- network Common Data Form.
- A self-describing, machine-independent data format.
- Probably the most common data format in the climate sciences.
- The names of netCDF files usually end with .nc.
- For further information see:

<http://www.unidata.ucar.edu/software/netcdf/>

# netCDF

- Load netCDF by entering the command:  
    `module load netcdf`
- Use `ncdump` to examine the contents of the sample atmosphere model output file, `stsc_spi62.nc`. Try commands such as:  
    `ncdump -h stsc_spi62.nc`  
    `ncdump -c stsc_spi62.nc`
- What can you see?



# Ferret

- A free visualisation and analysis package designed for visualising climatic data.
- For further information see:  
<http://ferret.pmel.noaa.gov/Ferret/>

## Using Ferret

- 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.

# Basic Ferret commands

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

# Basic Ferret commands

- 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 140E, 35S
<code>tsc[x=90e:180e,y=45s:0]</code>	Temperature within 90-180E,45-0S
<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

# Using Ferret

- 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[x=140e,y=35s,l=@ave]
```

```
list tsc[i=@ave,j=@ave,k=@ave,l=@ave]
```

```
show transform
```

# Using Ferret

- A sample ocean model output file, `com.spi62.00001.nc`, is also provided. Examine the contents of this file using `ncdump` and `Ferret`.

- Within `Ferret`, try commands such as:

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

- Try commands such as:

```
fill temp[i=@ave,l=@ave]  
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]
```

- Table 6.1 of the Users Guide will be useful here.

# Using Ferret

- The atmosphere data are output every pressure level in a different file, so we provide a script to combine them, in a single file:

```
comb_atm
```

- Look at the combined atmospheric outputs:

```
use su_exp01_all_atm.nc  
fill u[i=39,l=1]
```

```
use sv_exp01_all_atm.nc  
fill v[j=42,l=1]
```

# Using Ferret

- The ocean data are output every year in a different file, so we provide a script to combine them, in a single file:

```
comb_ocn
```

- Look at the combined atmospheric outputs:

```
use com.ct101.res_all_yr.nc  
fill/LEVELS=(-2,22,1)(inf) temp[i=@ave,l=@ave]
```

# Running a 50 years simulation

- The script `qsub_raijin_ct101` runs the coupled model for 50 years. It is the script that was used to run the control experiment.
- Using the `vi` command, examine this script.
- How does it differ from the script which runs the model for one day?

Hint: to compare two files, you can use `vimdiff`

```
vimdiff qsub_raijin_ct101 qsub_test_cpl
```

To exit, ESC then `:q!` twice (once for each file)



# Running a 50 years simulation

- Set up of the experiment part of the script

```
# Set name of run
set run = ctl01

# Set duration of run, in years
set DURATION = 50

# Set stack sizes
limit stacksize unlimited
setenv KMP_STACKSIZE 16M

setenv OMP_NUM_THREADS 10

# Create a directory for the outputs of this experiment
# if it already exists, delete the content
set TMP_DIR = /g/data/hh5/WS2019/$USER/$run/
if (-e $TMP_DIR) /bin/rm $TMP_DIR/*

# Copy the model executable to the run directory
cp $Mk3LHOME/core/bin/convert_averages $TMP_DIR

# Copy the control input file to the run directory
cp input_ctl01 $PBS_JOBFS/input
```

# Running a 50 years simulation

- The script that you just examined includes the following lines:

```
#PBS -l walltime=35:00:00
```

```
#PBS -l mem=2GB
```

```
#PBS -l ncpus=10
```

```
#PBS -l jobfs=2GB
```

- These request the resources needed to run the job.
- The job is expected to take up to 35 hours (walltime).
- The job will require up to 2 GB of memory (mem).
- We want to run on 10 cpus.
- We allocate 2GB of memory for the run directory
- When you design your own experiments, walltime is the only option that you might need to change. It takes ~24h for a 50 years run.

# Experiment setup

- To launch your experiment submit the script in the queue e.g.

```
qsub qsub_raijin_exp01
```

- To setup an experiment, you need to modify the input files taken by this script and the name of the experiment (run = ctl01)
- Remember that the three steps involved in running the model are:
  - create a run directory
  - copy model executable and input files to this directory
  - run the model

# Setup your own experiment

- The model requires three types of input files:
  - control file** configures the model for a particular simulation
  - restart files** initialise the model at the start of a simulation
  - auxiliary files** provide the boundary conditions during 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.

# Setup your own experiment

- Bottom boundary conditions:
  - topography
  - bathymetry
  - albedo
  - vegetation and soil types
- Radiative boundary conditions:
  - CO<sub>2</sub> transmission coefficients
  - ozone mixing ratios
- Applying a perturbation:
  - freshwater hosing mask

# Specific experiments

1. Magnitude of CO<sub>2</sub> effect on stormtracks
2. Reversibility of atmospheric CO<sub>2</sub>
3. Different epoch (mid-Holocene)
4. Difference between the melting of Greenland and Antarctic icesheets
5. Reflecting sunlight away through geoengineering
6. Effect of forests on rainfall