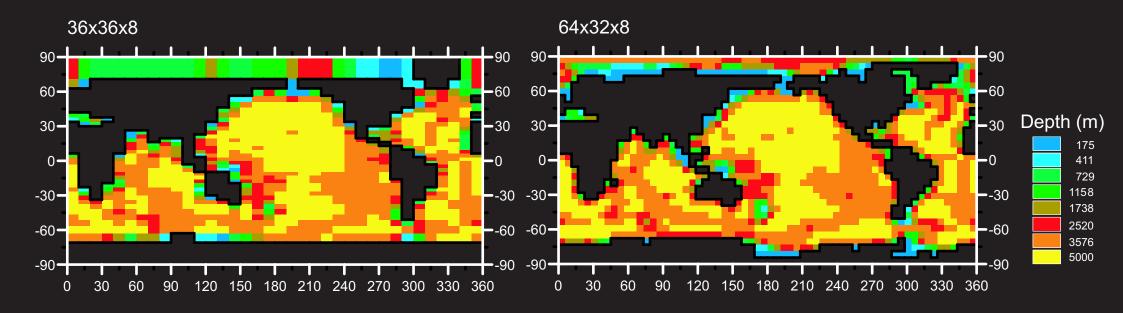
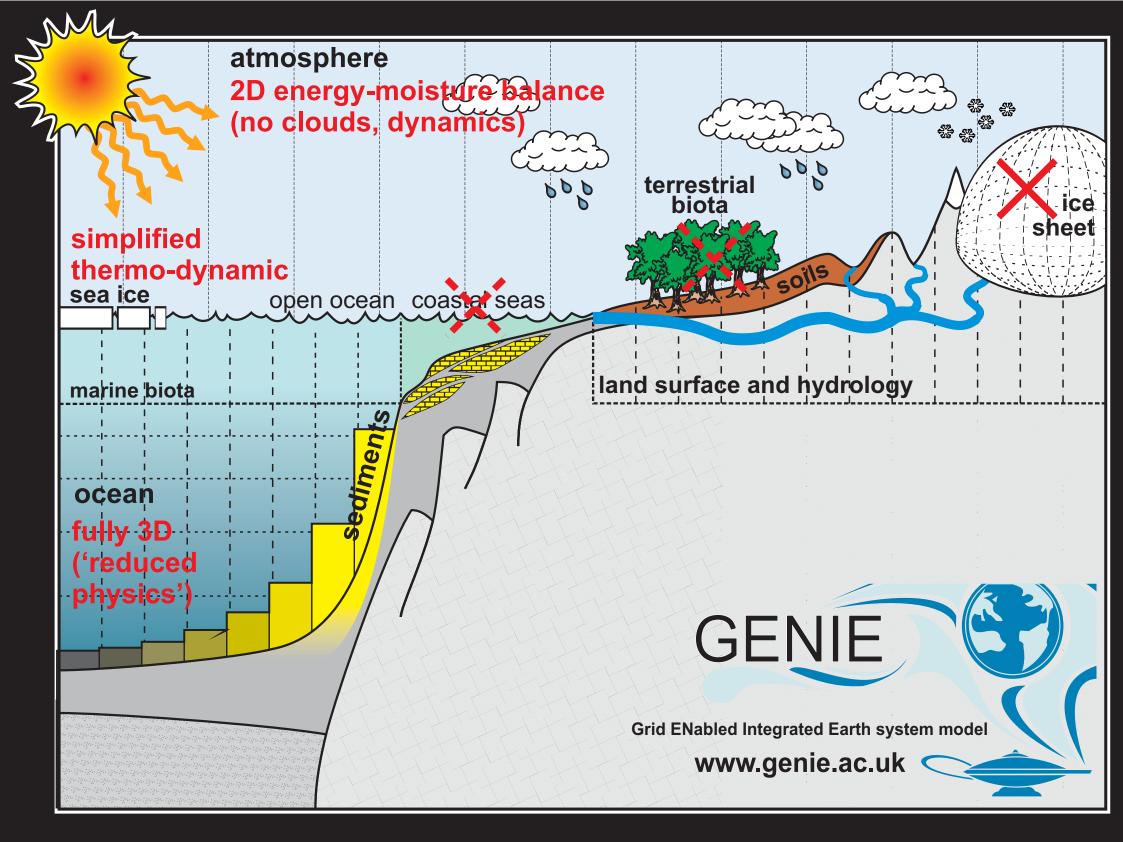


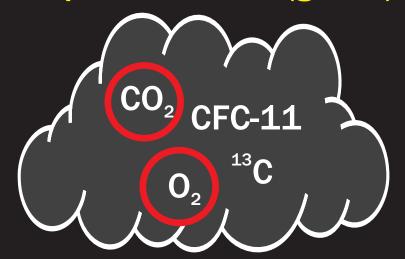
GENIE configuration

Model			gitude res.	`	gitude s res.	Vertical levels
Ocean GOLDSTEIN	\otimes	36 72 64 72	10° 5° 5.625° 10°	36 72 32 60	$sin(\pi/36)$ $sin(\pi/72)$ 5.625° 3°	8 16 8 8
Atmosphere 2-D EMBM	\otimes	36 72 64 72	10° 5° 5.625° 10°	36 72 32 60	$sin(\pi/36)$ $sin(\pi/72)$ 5.625° 3°	1 1 1
Atmosphere 3-D IGCM	\otimes	64	5.625°	32	5.625°	7

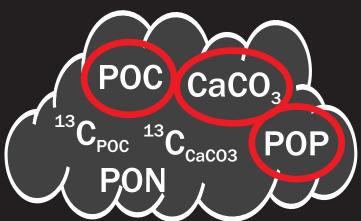




atmospheric tracers (gases)



biogeochemistry solid tracers (particulates)

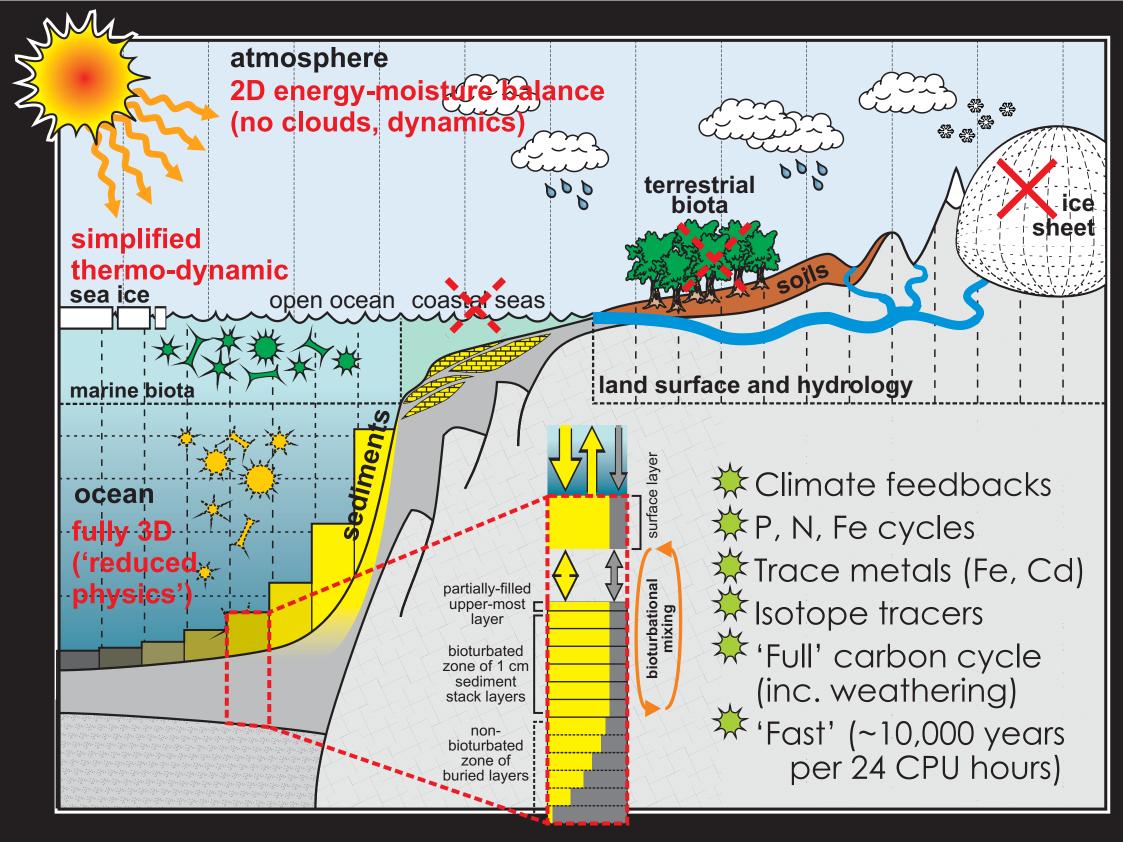


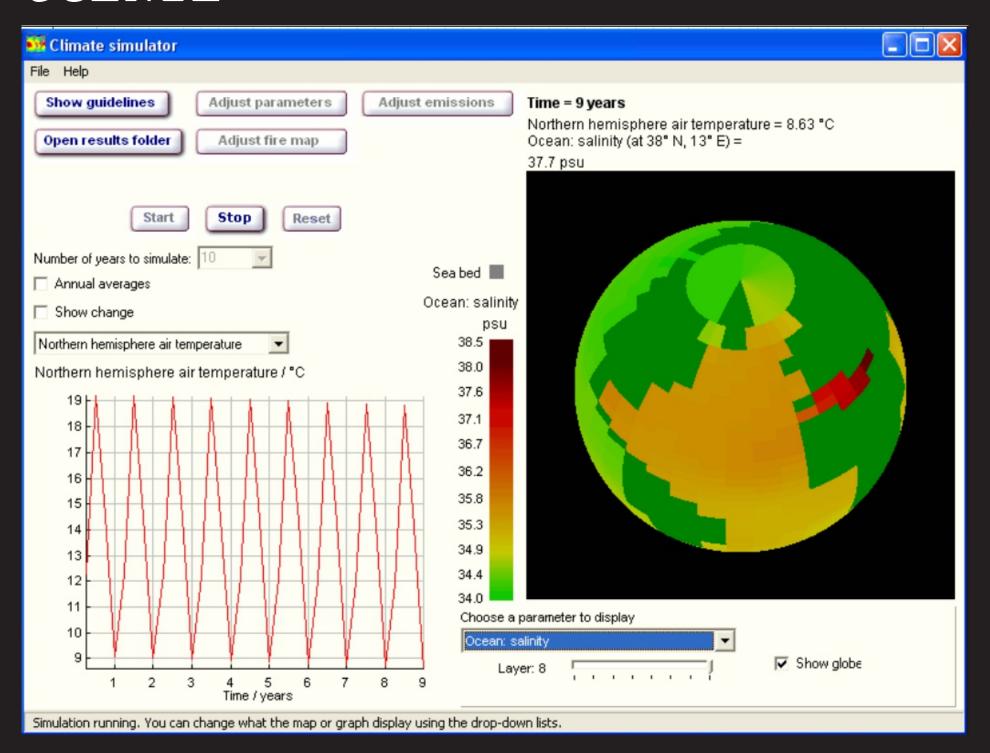
solubility coefficient Schmidt number PRE-DEFINED RELATIONSHIPS

e.g., Redfield ratios



dissolved tracers





! running GENIE

```
./old_rungenie.sh #1 #2 #3 #4 (#5)

parameter #1: 'base config' (e.g., cgenie_eb_go_gs_ac_bg)

parameter #2: 'user config' directory (e.g., LABS)

parameter #3: 'user config' file (e.g., exp1_modern)

parameter #4: experiment duration (e.g., 11)

parameter #5: restart experiment name and path (e.g., ~/genie_output/exp0_spinup)
```

! directory issues

cGENIE via runcgenie.sh and make cleanall are run from ~/genie/genie-main

You can change to ~/genie/genie-main from ~, either incrementally:

cd genie - main

or directly:

cd genie/genie-main
(or cd ~/genie/genie-main)
Return with cd on its own.

! directory viewing

Don't forget – the RH panel (the remote file system) in the SSH File Transfer Window does not automatically refresh the listing (and hence new experiment directories will be invisible until you refresh).

! make cleanall

This needs to be run only when the continental configuration, resolution of the ocean model, or number of 'tracers' in the ocean have been changed. (You will be advised when this is essential.) BUT NOTE: If you have carried out a make cleanall, BEFORE you can submit an experiment to the cluster the model must be recompiled – i.e., simply run the model interactively for a few years. Note that it is 'good practice' that when creating a completely new sort of model experiment (i.e., not just simply changing again a single parameter value), that you try running it for a few years interactively, as this will tell you immediately if there is something incorrectly set up in the experiment.

! experiment run length and results saving By default: cGENIE saves time-slice (netCDF format) results at times centred on years:

0.5 1.5 4.5 9.5 19.5 49.5 99.5 199.5 ... (a similar pseudo-log sequence is used for time-series saving) Hence, to capture the final state of the model at the end of an

experiment, the experiment length has to be one of:

```
1 2 5 10 20 50 100 200 ...
```

as the time-slices are annual averages by default (and saved at the mid-point of the average).

The default time-points for time-slice saving can be changed (via a simple text file), found in ~/cgenie/genie-biogem/data/input and called: save_timeslice.dat

(for time-series saving, the file is called: save_timeseries.dat)

! editing experiments

Simply: edit the user config text file, and save!

! creating new experiments

Is as simple as copying a pre-existing (and presumably similar) user config file and giving it a different name.

Either:

```
cp exp1file exp2file
cp snowball x10co2 snowball x20co2
```

!(then edit it)or, all within Windoz:

Using the file transfer client:

- (1) transfer a user config file to your local PC (or remote UoB filespace),
- (2) rename it
- (3) transfer it back(then edit it)