The background of the image is a vast, calm body of water under a sky filled with large, wispy clouds. The colors transition from deep orange and red at the horizon to a lighter yellow and white higher up. In the upper left foreground, a dark, silhouetted shape, possibly a rock or a small boat, is visible against the bright sky. The overall atmosphere is serene and suggests a polar or arctic environment.

# **COAWST configuration guide**

**Third Edition**

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The *model2roms* toolbox (<https://github.com/trondkr/model2roms>) was created by Trond Kristiansen ([me@trondkristiansen.com](mailto:me@trondkristiansen.com)) and is licensed under a *MIT License*.



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**Third edition**

September 2020





## Reviewers

The authors would like to thank the following peers for reviewing the **COAWST configuration Guide - First Edition**:

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Thank you.





## Author's note

This guide is designed to assist new users to use the Coupled Ocean-Atmosphere-Wave-Sediment Transport System (COAWST). The main idea behind this guide is to teach the necessary steps to use COAWST, beginning with its installation, then a simulation of a test case and the configuration of a project. To achieve this goal, we use several programming languages, such as Fortran, Python and MATLAB. In the future we intend to adapt all scripts to a free programming language.

When we started writing this guide, we wanted to pass on our experience of using a numerical modeling system that is considered the state of the art in our field, throughout reading, understanding how it works and how to use the COAWST, allying theory with practice.

However, a major difficulty in this process was the generation of the boundary and initial conditions of the oceanic model, the Regional Ocean Modeling System (ROMS), which rely on paid software. To get around this problem we chose to work with the *model2roms* toolbox package. This set of routines was developed in Python and Fortran language by Trond Kristiansen (<http://www.trondkristiansen.com>).

We emphasize that in some chapters, the reader will find how to use the COAWST in a cluster that is available for use by the Ocean and Atmosphere Studies Laboratory (LOA) of the National Institute for Space Research (INPE). This is a system of high performance computing that allows parallel numerical operations. In this case, the guide may serve only as inspiration and the reader's prior knowledge in applying COAWST in his own cluster system will be worthwhile.

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UnderConstruction

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We wish you a good reading and success in your research.

The authors.



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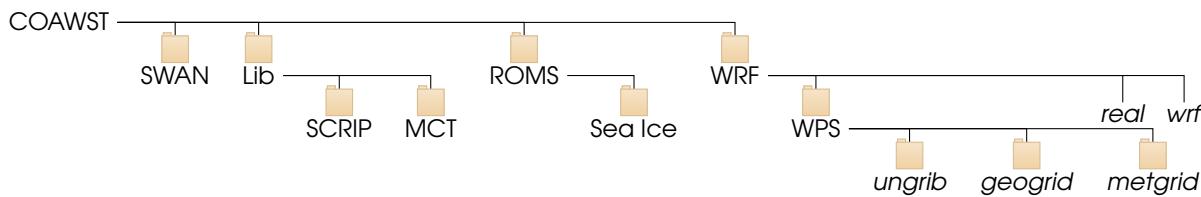




# 1. Introduction

This guide provides a brief introduction to the models that make up the Coupled Ocean-Atmosphere-Wave-Sediment Transport Modeling System (COAWST), as well as the additional user settings required for specific projects. As shown in Figure 1.1, COAWST uses several models and programs that will be presented below:

- **COAWST**: Core of the coupled numerical modeling system. More information in the section 1.7;
- **ROMS**: The hydrodynamic model. More information in the Section 1.1;
- **Sea Ice**: The sea ice model, coupled to ROMS. In this guide we use the Budgell's Sea Ice Model. More information in the Section 1.2;
- **WRF**: The atmospheric model. More information in the Section 1.3;
- **wrf**: Executable program to start the WRF atmospheric simulation. More information in the section 4.4;
- **real**: Executable program to generate the initial condition and the boundary forces of the WRF. More information in the Section 4.3.5;
- **WPS**: Package with three programs to generate the files to be used in *real*. More information in the Section 4.2;
- **geogrid**: Program to, mainly, generate the WRF grid domain. More information in the section 4.3.1;
- **ungrib**: Briefly, the program that extracts the data in *GRIB* format. More information in the Section 4.3.3;
- **metgrid**: Briefly, the program that interpolates the data generated by *ungrib*. More information in the Section 4.3.4;
- **SWAN**: The wave model. More information in the Section 1.4;
- **MCT**: The set of codes that couples the previous models. More information in the Section 1.5;
- **SCRIP**: Package that interpolates and remaps the model's grids to allow the models to be coupled. More information in the section 1.6.



**Figure 1.1.** COAWST folder structure

When writing this guide, we chose to use Python as a programming language for some steps. It is a programming language designed with a philosophy of emphasizing the importance of the programmer's effort over the computational effort and prioritizes code readability over speed or expressiveness. The language has a wide and active community, which facilitates the search information by the user, as there is an extensive collection of libraries and documents on the internet.

Python Brazil (<http://python.org.br>) offers great help for beginners, providing introductions about the language and also a guide to use Python for scientific purposes.

We also use MATLAB as a language. It is a paid, but high-performance interactive software focused on the numerical calculation. The MATLAB Answers website is a platform created to offer help to the users about the language. The site is available at <https://www.mathworks.com/matlabcentral/answers/index>.

## 1.1 Regional Ocean Modeling System

The Regional Ocean Modeling System (ROMS; [Shchepetkin & McWilliams, 2005](#)) is a three-dimensional ocean model with free surface, vertical sigma coordinate with sigma vertical coordinates (that follow the terrain) and solve primitive equations. The model uses the Reynolds average and the finite difference method to solve the Navier-Stokes equations using hydrostatic approximations and Boussinesq ( cite Haidvogel2008).

The hydrostatic equations of momentum use a split-explicit time-step scheme, where the barotropic and baroclinic modes are solved separately, in different finite numbers of steps of time, to solve the free surface equations and it is integrated vertically. This structure separate time-steps frames maintains the volume conservation and consistency preservation that are necessary for the tracers ([Haidvogel et al., 2008](#); [Shchepetkin & McWilliams, 2005](#)).

The model solves the horizontally equations through orthogonal curvilinear coordinates of the Arakawa-C grid type ([Arakawa & Lamb, 1977](#)). Vertically, the coordinates follow the features of the terrain and allow you to adjust the resolution along the water column. To guarantee the conservation of momentum, the grid uses second order finite differences ([Haidvogel et al., 2008](#)).

ROMS is a model that has free code and its development has the contribution of the user community. Currently, the version used in COAWST is managed by Dr. Hernan Arango of Rutgers University. To access the model code, is necessary to register on the ROMS website (<https://www.myroms.org/>) The site has a extremely useful and very active forum to discuss about questions and suggestions. You can access here: (<https://www.myroms.org/forum>)

We recommend to read the ROMS Technical Manual, written by [Hedström \(2018\)](#). This manual has several information about the equations and algorithms of the model and examples of test cases.

## 1.2 Budgell's Sea Ice Model

The Sea Ice Model, proposed by [Budgell \(2005\)](#), has the same time and grid steps as the ROMS model and shares the same parallel encoding structure for use with Message Passing Import (MPI). Thus, allows dynamic and thermodynamic modeling where sea ice predominates, such as at high latitudes.

The main attributes of the model, according to [Hedström \(2018\)](#), are:

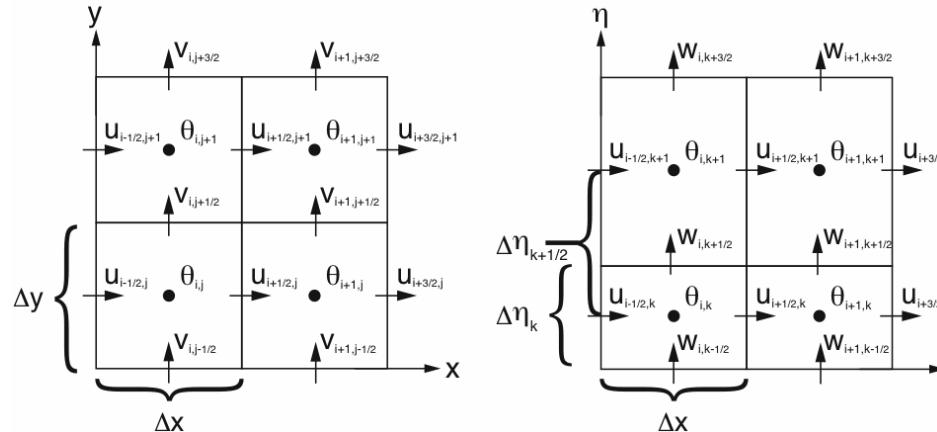
- [Hunke & Dukowicz \(1997\)](#) and [Hunke \(2001\)](#) elastic-viscous-plastic dynamics;
- [Mellor & Kanta \(1989\)](#) thermodynamics;
- Orthogonal-curvilinear coordinates;
- [Arakawa & Lamb \(1977\)](#) grid;
- [Smolarkiewicz & Grabowski \(1990\)](#) advection of tracers;
- [Lemieux et al. \(2015\)](#) landfast ice parameterization.

## 1.3 Weather Research & Forecasting Model

The Weather Research and Forecasting (WRF; [Skamarock et al., 2008](#)) is a model developed by the National Centers for Environmental Prediction (NCEP), the National Center for Atmospheric Research (NCAR) and research groups from different universities.

To integrate the governing equations over time, the Advanced Research WRF (ARW) uses low frequency modes that are integrated using the third-order Runge-Kutta scheme, and the integrated acoustic and gravity (high frequency) modes with a lower time step. By this way, the numerical stability is maintained through a “forward-backward” scheme for acoustic modes that propagate horizontally and an implicit scheme for acoustic modes for vertical propagation and buoyancy oscillations ([Skamarock et al., 2008](#)).

The WRF model uses an Arakawa-C type grid ([Arakawa & Lamb, 1977](#)), where normal speeds are staggered halfway through the grid of thermodynamic variables, as shown in the schematic representation illustrated in Figure 1.2.



**Figure 1.2.** Horizontal and vertical grid of Weather Research and Forecast (WRF) using the Arakawa-C grid. The horizontal and vertical components of velocity ( $u$ ,  $v$  and  $w$ ) are positioned along the faces of the grids and the thermodynamic variables ( $\theta$ ) are positioned in the center of each grid.

Author: [Skamarock et al. \(2008\)](#).

It is important to note that the WRF, without coupling with other models, simulates the surface roughness based on the ratio of roughness to wind shear proposed by Charnock (1955), as exemplified in the following equation 1.1:

$$Z_0 = Z_{ch} \frac{u_*^2}{g} \quad (1.1)$$

Where  $Z_0$  é a roughness,  $Z_{ch}$  is the Charnock parameter (a dimensionless value of 0,018),  $u_*$  the frictional speed (m/s) and  $g$  the gravity acceleration(9,81 m/s<sup>2</sup>).

The WRF is available at: [http://www2.mmm.ucar.edu/wrf/users/download/get\\_source.html](http://www2.mmm.ucar.edu/wrf/users/download/get_source.html).

## 1.4 Simulating Waves Nearshore

The Simulating Waves Nearshore (SWAN; Booij et al., 1996, 1999) is a third generation model, designed to simulate coastal regions with shallow waters and local currents. The model is widely used in the numerical forecast of the Simulating Waves Nearshore (SWAN; Booij et al., 1996, 1999) is a third generation model, designed to compute in coastal regions with shallow waters and local currents. The model is widely used in the numerical forecast of waves in coastal regions, estuaries, channels and others, being able to use fields of wind, bathymetry and currents provided by other models (Booij et al., 1996, 1999).

Silva (2013) and Booij et al. (1996; Booij et al., 1999) list the main characteristics of the SWAN:

- Wave propagation in time and space, shoaling, refraction due to current and depth, frequency shifting due to currents and non-stationary depth;
- Wave generation by wind;
- Whitecapping, bottom friction and depth-induced breaking;
- Dissipation due to aquatic vegetation, turbulent flow and viscous fluid mud;
- Wave-induced set-up;
- Propagation from laboratory up to global scales;
- Transmission through and reflection (specular and diffuse) against obstacles;
- Diffraction.

More features can be found in SWAN website, available at: <http://swanmodel.sourceforge.net/>.

## 1.5 Model Coupling Toolkit

The Model Coupling Toolkip (MCT; Jacob et al., 2005; Larson et al., 2005; Warner et al., 2008) is a set of open-source scripts, written in Fortran90 that allow the transmission and transformation of the different data necessary for model coupling. During initialization, model domains are broken down into segments that are distributed between processors, allowing models to be coupled also in parallel.

According to the MCT website, (<http://www.mcs.anl.gov/research/projects/mct/>), the toolkit provides the following core coupling services:

- A component model registry;
- Domain decomposition descriptors;
- A time averaging and accumulation buffer datatype;
- A general spatial grid representation capable of supporting unstructured grids;
- Parallel tools for intergrid interpolation;
- Tools for merging data from multiple components for use by another component;
- A programming model similar to that of the Message Passing Interface.

## 1.6 Spherical Coordinate Remapping Interpolation Package

The Spherical Coordinate Remapping Interpolation Package (SCRIP; [Jones, 1998, 1999](#)) is freely available for download at <https://github.com/SCRIP-Project/SCRIP>. The package is distributed together with COAWST modeling system. This package is used for projects that use more than one model and with different grids (with different spatial resolutions). SCRIP will generate the interpolation weights that will be used to remap the data between the different grids of the different models.

In COAWST, SCRIP was modified to generate a single file (in NetCDF format) that contains the weights based in each model grid.

## 1.7 Coupled-Ocean-Atmosphere-Wave-Sediment Transport Modeling System

The COAWST [Warner et al., 2010, 2008](#)) uses the WRF as the atmospheric model, the ROMS as the hydrodynamic model, the SWAN as the wave model and the sediment transport model Community Sediment Transport Modeling Project (CSTM; [Warner et al., 2008](#)), each one coupled by the MCT ([Warner et al., 2010, 2008](#)). The frequency with which this information is exchanged between the different models is adjusted by the user.

The coupling between the models allows the different physical processes that occur in both oceanic and atmospheric environments to be identified and analyzed with greater accuracy when compared to simulations without active coupling. ([Miller et al., 2018; Pullen et al., 2018](#)).

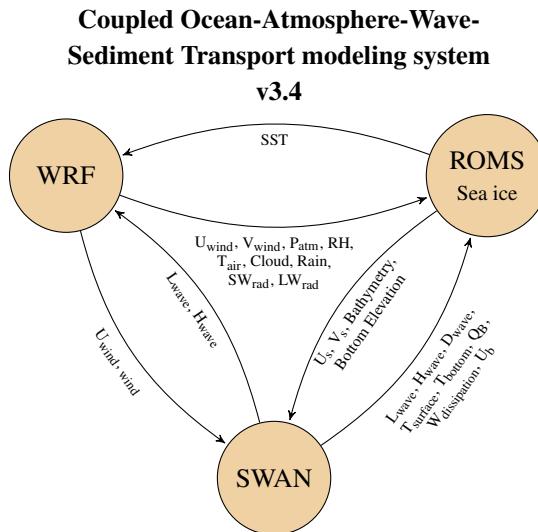
**WARNING**

This guide does not use CSTM. In case you are interested, there is a study on the transfer of sediments during Hurricane Isabel (2003) by [Warner et al. \(2010\)](#).

As shown in Figure 1.3, the informations exchanged between models are:

- WRF -> ROMS: surface shear and liquid heat fluxes (calculated in ROMS from the components of latent and sensitive heat fluxes) shortwave and longwave radiation, atmospheric pressure, relative humidity, air temperature, clouds, precipitation and wind components ;
- ROMS -> WRF: sea surface temperature;
- SWAN -> ROMS: surface and bottom wave direction, height, length, period, energy dissipation and lower orbital speed;
- ROMS -> SWAN: bathymetry, surface elevation, height of the sea and average currents in depth;

- SWAN -> WRF: roughness of the sea surface (calculated in WRF from the significant wave height, length and period);
- WRF -> SWAN: wind at 10m.



**Figure 1.3.** Diagram about the exchange of information between the main models that make up the COAWST modeling system. (Warner et al., 2008)

The Woods Hole Coastal and Marine Science Center page provides an experimental presentation on real time of COAWST (Sea Surface Temperature, Sea Surface Height, Significant Height Wave, Current and Wind Vectors and Sediment Dispersion) in the eastern United States and Gulf of Mexico. The content is available at: <https://woodshole.er.usgs.gov/project-pages/cccp/public/COAWST.htm>.

## 1.8 Requirements to use this guide

To compile COAWST, this guide uses a cluster with parallel communication and a nonparallel Linux OS with Gfortran. If you choose to compile in on a computer with different specifications, use the original COAWST user manual as a reference, available in the modeling system repository. In this case, see how to download COAWST in the [2.6](#) section.

We use **Ubuntu 19.04 LTS**. It is important to maintain the same systems operational, as other version may conflict with some libraries used by the *model2roms* toolkit or even the COAWST itself.

**WARNING** This guide uses COAWST v. 3.6.



## 2. COAWST in a parallel cluster

### 2.1 About our cluster Kerana

The CRAY XE machine, also called Kerana, is a cluster with massively parallel architecture, with 84 processing nodes and 2688 cores and is located at CPTEC/INPE, in Cachoeira Paulista, São Paulo. For having the ability to parallelize operations through the MPI interface, the cluster is ideal for using numerical models with high spatial resolution and temporal.

### 2.2 Signing up a user account

To start the process of applying for a user account in Kerana, it is necessary that the computer in the INPE's premises have a fixed IP. Your advisor or supervisor is required to send an email to *Helpdesk* ([helpdesk.cptec@inpe.br](mailto:helpdesk.cptec@inpe.br)) informing the MAC address, hostname and reason for the request. If the computer already has a fixed IP assigned, it also need to be informed to be changed.

With the fixed IP configured, contact the *Helpdesk* ([helpdesk.cptec@inpe.br](mailto:helpdesk.cptec@inpe.br)) to request a form to use Kerana.

### 2.3 Signing in to the Kerana cluster

Access will be done entirely through the computer terminal. You will need two primary commands: *ssh* for accessing and manipulating files and folders and *sftp* to download and upload files.

To access and modify files and folders, type in the terminal, replacing *name.surname* with the user provided by *Helpdesk*:

**WARNING**

From now on, whenever the guide shows the user as *name.surname*, change to your username provided by the Helpdesk.

```
1 ssh -Y name.surname@acesso-hpc.cptec.inpe.br -p 2000
```

To download and/or upload, use:

```
1 sftp -P2000 name.surname@acesso-hpc.cptec.inpe.br
```

**WARNING**

It is not possible to download and upload multiple files at the same time using *sftp*. One tip is to compress into a single *tar* file and then unzip them.

To add files from your computer to Kerana:

```
1 put file.tar.gz
```

To extract Kerana files to your computer:

```
1 get file.tar.gz
```

## 2.4 File repository

Certain files are needed in each user's area to facilitate the use of the cluster. You will find them in the directory:

```
1 /scratch/adriano.sutil/repositorio/
```

To copy the files to your area, type:

```
1 cp -r /scratch/adriano.sutil/repositorio /scratch/name.surname
```

**WARNING**

From now on this guide will use the files that are inside this repository, so it is essential that they are in your area.

## 2.5 Kerana environment

It is necessary to activate some modules in the cluster to compile COAWST. In this case, open the file `.bashrc` which is located at the root directory of your user.

```
1 vim .bashrc
```

Add the following commands at the end of the file, changing only `name.surname` to your username:

```
1 module load java
2 module load netcdf
3
4 export PATH=/scratch/name.surname/repositorio/Softs/nedit/5.5:$PATH
5 export PATH=/scratch/name.surname/repositorio/Softs/bin:$PATH
6
7 export PHDF5=${HDF_DIR}
8 export WRFIO_NCD_LARGE_FILE_SUPPORT=1
9
10 export PATH=/home/luciano.pezzi/local/bin:$PATH
11 export JASPERINC=/home/luciano.pezzi/local/include
12 export JASPERLIB=/home/luciano.pezzi/local/lib
13 export LD_LIBRARY_PATH=/home/luciano.pezzi/local/lib:$LD_LIBRARY_PATH
```

Save and type in the terminal:

```
1 source .bashrc
```

## 2.6 Downloading COAWST

**WARNING**

COAWST v3.6 is already in the repository within the Kerana cluster, as discussed in Section 2.4.

To download COAWST, send an email to Dr. John Warner ([jcwarner@usgs.gov](mailto:jcwarner@usgs.gov)), one of the heads behind the coupled regional modeling system.

After access is granted, with the user credentials and password provided by Dr. John Warner, type in the command the command below, changing the `textit myusrname` to your username.

```
1 svn checkout --username myusrname https://coawstmodel.sourcerepo.com/coawstmodel/COAWST
```

Add the COAWST folder to your Kerana desktop using `sftp`, as in the Section 2.3.

## 2.7 Automating the compilation of COAWST in Kerana

**WARNING** This guide uses COAWST v3.6.

To speed up the process, it is possible to automate some compilation steps. Enter the directory:

```
1 cd /scratch/name.surname/COAWST/WRF/arch
```

Open the file *Config.pl*:

```
1 nedit Config.pl
```

Look for the lines:

```
1 printf "\nEnter selection [%d-%d] : ",1,$opt ;
2 $response = <STDIN> ;
```

And replace <STDIN> to 42, as in the example below:

```
1 printf "\nEnter selection [%d-%d] : ",1,$opt ;
2 $response = 42 ;
```

Open the file *configure.defaults*:

```
1 nedit configure.defaults
```

Look for the *TRADEFLAG*, option, which is on line 1262, approximately:

```
1 TRADEFLAG = CONFIGURE_TRADEFLAG
```

And modify to:

```
1 TRADEFLAG = -traditional
```

These modifications will select the configurations of the Kerana cluster (*CRAY CCE (ftn/gcc)*: *Cray XE* and *XC (dmpar)*), among those available to use COAWST, as shown in Figure 2.1.

1. (serial)	2. (smpar)	3. (dmpar)	4. (dm+sm)	PGI (pgf90/gcc)
5. (serial)	6. (smpar)	7. (dmpar)	8. (dm+sm)	PGI (pgf90/pgcc): SGI MPT
9. (serial)	10. (smpar)	11. (dmpar)	12. (dm+sm)	PGI (pgf90/gcc): PGI accelerator
13. (serial)	14. (smpar)	15. (dmpar)	16. (dm+sm)	INTEL (ifort/icc)
18. (serial)	19. (smpar)	20. (dmpar)	21. (dm+sm)	INTEL (ifort/icc): Xeon Phi (MIC architecture)
22. (serial)	23. (smpar)	24. (dmpar)	25. (dm+sm)	INTEL (ifort/icc): Xeon (SNB with AVX mods)
26. (serial)	27. (smpar)	28. (dmpar)	29. (dm+sm)	INTEL (ifort/icc): SGI MPT
30. (serial)		31. (dmpar)		INTEL (ifort/icc): IBM POE
32. (serial)	33. (smpar)	34. (dmpar)	35. (dm+sm)	PATHSCALE (pathf90/pathcc)
36. (serial)	37. (smpar)	38. (dmpar)	39. (dm+sm)	GNU (gfortran/gcc)
40. (serial)	41. (smpar)	42. (dmpar)	43. (dm+sm)	IBM (xlf90_r/c_r_r)
44. (serial)	45. (smpar)	46. (dmpar)	47. (dm+sm)	PGI (ftn/gcc): Cray XC CLE
48. (serial)	49. (smpar)	50. (dmpar)	51. (dm+sm)	CRAY CCE (ftn/gcc): Cray XE and XC
52. (serial)	53. (smpar)	54. (dmpar)	55. (dm+sm)	INTEL (ftn/icc): Cray XC
56. (serial)	57. (smpar)	58. (dmpar)	59. (dm+sm)	PGI (pgf90/pgcc)
60. (serial)	61. (smpar)	62. (dmpar)	63. (dm+sm)	PGI (pgf90/gcc): -f90=pgf90
				PGI (pgf90/pgcc): -f90=pgf90

**Figure 2.1.** Computational options available for selection when compiling COAWST.

Now, open the *Config\_new.pl* file and then look for the following lines:

```
1 printf "Compile for nesting? (1=basic, 2=preset moves, 3=vortex following) [default 1]: " ;
2 }
3 $response = <STDIN> ;
```

And change <STDIN> to the basic WRF atmospheric model nesting mode, as in the following example:

```
1 printf "Compile for nesting? (1=basic, 2=preset moves, 3=vortex following) [default 1]: " ;
2 }
3 $response = 1 ;
```

## 2.8 Compiling the MCT

**WARNING** This guide uses COAWST v 3.6.

Each new user must compile the MCT before compiling COAWST. The first step is to use the *setup\_pgi.sh* file. This file was copied from the previous repository root, it is essential to change the directories contained therein. So open the file:

```
1 nedit setup_pgi.sh
```

If necessary, change the directories according to the name of your COAWST folder and execute the file to load the necessary modules:

```
1 source setup_pgi.sh
```

The libraries will be activated, as shown in Figure 2.2:

```
Currently Loaded Modulefiles:
 1) modules/3.2.6.7
 2) nodestat/2.2-1.0400.29866.4.3.gem
 3) sdb/1.0-1.0400.31073.9.3.gem
 4) MySQL/5.0.64-1.0000.4667.20.1
 5) lustre-cray_gem_s/1.8.4_2.6.32.45_0.3.2_1.0400.6336.8.1-1.0400.30879.1.81
 6) udreg/2.3.1-1.0400.3911.5.13.gem
 7) ugni/2.3-1.0400.4127.5.20.gem
 8) gni-headers/2.1-1.0400.4156.6.1.gem
 9) dmapp/3.2.1-1.0400.3965.10.63.gem
10) xpmem/0.1-2.0400.30792.5.6.gem
11) hss-llm/6.0.0
12) Base-opts/1.0.2-1.0400.29823.8.5.gem
13) xtpe-network-gemini
14) cce/8.6
15) totalview-support/1.1.3
16) xt-totalview/8.10.0
17) acml/5.1.0
18) xt-libsci/11.1.01
19) pmi/3.0.1-1.0000.9101.2.26.gem
20) rca/1.0.0-2.0400.30002.5.75.gem
21) xt-asyncpe/5.14
22) atp/1.5.1
23) PrgEnv-cray/4.0.36
24) pbs/10.4.0.101257
25) xt-mpich2/5.5.4
26) xtpe-interlagos
27) java/jdk1.7.0_07
28) grads/2.0.a8

Currently Loaded Modulefiles:
 1) modules/3.2.6.7
 2) nodestat/2.2-1.0400.29866.4.3.gem
 3) sdb/1.0-1.0400.31073.9.3.gem
 4) MySQL/5.0.64-1.0000.4667.20.1
 5) lustre-cray_gem_s/1.8.4_2.6.32.45_0.3.2_1.0400.6336.8.1-1.0400.30879.1.81
 6) udreg/2.3.1-1.0400.3911.5.13.gem
 7) ugni/2.3-1.0400.4127.5.20.gem
 8) gni-headers/2.1-1.0400.4156.6.1.gem
 9) dmapp/3.2.1-1.0400.3965.10.63.gem
10) xpmem/0.1-2.0400.30792.5.6.gem
11) hss-llm/6.0.0
12) Base-opts/1.0.2-1.0400.29823.8.5.gem
13) xtpe-network-gemini
14) PrgEnv-pgi/4.0.36
15) xt-mpich2/5.5.4
16) atp/1.5.1
17) xt-asyncpe/5.14
18) pmi/3.0.1-1.0000.9101.2.26.gem
19) xt-libsci/11.1.01
20) xt-totalview/8.10.0
21) totalview-support/1.1.3
22) pgi/12.8.0
23) pbs/10.4.0.101257
24) xtpe-interlagos
25) java/jdk1.7.0_07
26) grads/2.0.a8
27) hdf5-parallel/1.8.8
28) netcdf-hdf5parallel/4.2.0
29) libfast/1.0.9

adriano.util@login1:~/COAWST_WR> █
```

**Figure 2.2.** Modules activated in the cluster with the file *setup\_pgi.sh* with user adriano.util.

Enter the MCT folder directory:

```
1 cd /home/name.surname/COAWST/Lib/MCT
```

Open the file *Makefile.conf*:

```
1 nedit Makefile.conf
```

And modify the file as follows:

**WARNING** Remember to change the *name.surname*!

```

1 FC          = ftn
2 FCFLAGS     = -O2
3 F90FLAGS    =
4 REAL8       = -r8
5 ENDIAN      = -Mbyteswapio
6 INCFLAG     = -I
7 INCPATH     =
8 MPILIBS     =
9 DEFS        = -DSYSLINUX -DCPRPGI
10 FPP         = cpp
11 FPPFLAGS   = -P -C -N -traditional
12 CC          = cc
13 ALLCFLAGS   = -DFORTTRAN_UNDERSCORE_ -DSYSLINUX -DCPRPGI -O
14 COMPILER_ROOT =
15 BABELROOT   =
16 PYTHON      =
17 PYTHONOPTS  =
18 FORT_SIZE   =
19 CRULE       = .c.o
20 90RULE     = .F90.o
21 F90RULECPP = .F90RULECPP
22 INSTALL     = /home/name.surname/COAWST/Lib/MCT/install-sh -c
23 MKINSTALLDIRS = /home/name.surname/COAWST/Lib/MCT/mkinstalldirs
24 abs_top_builddir= /home/name.surname/COAWST/Lib/MCT/
25 MCTPATH     = /home/name.surname/COAWST/Lib/MCT/mct
26 MPEUPATH    = /home/name.surname/COAWST/Lib/MCT/mpeu
27 EXAMPLEPATH = /home/name.surname/COAWST/Lib/MCT/examples
28 MPISERPATH =
29 libdir      = /home/name.surname/COAWST/Lib/MCT/pgi/lib
30 includedir  = /home/name.surname/COAWST/Lib/MCT/pgi/include
31 AR          = ar cq
32 RM          = rm -f

```

Install the MCT by typing the following commands:

```

1 make
2 make install

```

Observe the messages that appear in the terminal and look for errors. If not, the MCT was successfully compiled.

## 2.9 Compiling the Sandy test case

There are some test cases within COAWST to be compiled and worked on. In this case we'll compile Hurricane Sandy's project, which couples and nests WRF, ROMS and SWAN. First, it is necessary to know

the structure of COAWST files and folders.

The typical COAWST directory structure is exemplified in Figure 2.3. Mainly, we will use the folders *Projects* e *Work* to work on.

```
drwxr-xr-x 20 adriano.sutil users 4096 Ago 3 13:04 .
drwxrwx--- 10 adriano.sutil users 4096 Ago 4 13:59 ..
drwxr-xr-x 6 adriano.sutil users 4096 Ago 3 15:25 ARWpost
drwxr-xr-x 2 adriano.sutil users 36864 Jul 26 16:28 Build
-rw-rxr-xr-x 1 adriano.sutil users 18212 Jul 14 12:37 coawst.bash
-rw-rxr-xr-x 1 adriano.sutil users 61866264 Jul 26 16:28 coawstM
-rw-r--r-- 1 adriano.sutil users 1163942 Jul 26 16:28 coawst.pgi.wrs
-rw-rxr-xr-x 1 adriano.sutil users 2126848 Jul 14 10:30 COAWST_User_Manual.doc
drwxr-xr-x 2 adriano.sutil users 4096 Jul 14 10:30 Compilers
drwxr-xr-x 3 adriano.sutil users 4096 Jul 14 10:30 Data
-rw-rxr-xr-x 1 adriano.sutil users 120 Jul 14 10:24 ..DS_Store
drwxr-xr-x 1 adriano.sutil users 12292 Jul 14 10:24 .DS_Store
-rw-rxr-xr-x 1 adriano.sutil users 261 Jul 14 10:30 GENPARM.TBL
drwxr-xr-x 8 adriano.sutil users 4096 Jul 14 10:30 InWave
-rw-rxr-xr-x 1 adriano.sutil users 29820 Jul 14 10:30 LANDUSE.TBL
drwxr-xr-x 5 adriano.sutil users 4096 Jul 14 10:31 Lib
-rw-rxr-xr-x 1 adriano.sutil users 23983 Jul 14 10:32 makefile
drwxr-xr-x 2 adriano.sutil users 4096 Jul 14 10:32 Master
drwxr-xr-x 3 adriano.sutil users 4096 Jul 26 15:37 Projects
-rw-rxr-xr-x 1 adriano.sutil users 3018 Jul 14 10:42 README.txt
drwxr-xr-x 2 adriano.sutil users 4096 Jul 14 10:32 REFDEF
drwxr-xr-x 16 adriano.sutil users 4096 Jul 14 10:33 ROMS
-rw-rxr-xr-x 1 adriano.sutil users 749248 Jul 14 10:33 RRTM_DATA
-rw-rxr-xr-x 1 adriano.sutil users 1498368 Jul 14 10:33 RRTM_DATA_DBL
-rw-rxr-xr-x 1 adriano.sutil users 500 Jul 14 10:33 run_coawst
drwxr-xr-x 3 adriano.sutil users 4096 Jul 14 10:33 SeaIce
-rw-rxr-xr-x 1 adriano.sutil users 739 Jul 14 10:40 setup_cray.sh
-rw-rxr-xr-x 1 adriano.sutil users 748 Jul 14 10:40 setup_pgi.sh
-rw-rxr-xr-x 1 adriano.sutil users 4419 Jul 14 10:33 SOILPARM.TBL
drwxr-xr-x 4 adriano.sutil users 4096 Jul 14 10:30 .SVN
drwxr-xr-x 4 adriano.sutil users 4096 Jul 14 10:33 SWAN
drwxr-xr-x 3 adriano.sutil users 4096 Jul 14 10:34 Tools
-rw-rxr-xr-x 1 adriano.sutil users 11190 Jul 14 10:34 URBPARM.TBL
drwxr-xr-x 5 adriano.sutil users 4096 Jul 14 10:34 User
-rw-rxr-xr-x 1 adriano.sutil users 22717 Jul 14 10:34 VEGPARM.TBL
drwxr-xr-x 3 adriano.sutil users 4096 Jul 26 15:38 Work
drwxr-xr-x 7 adriano.sutil users 20480 Jul 26 17:29 WPS
drwxr-xr-x 17 adriano.sutil users 4096 Jul 26 15:50 WRF
adriano.sutil@login1:~/COAWST_V3.2>
```

**Figure 2.3.** Representation of the main COAWST folder and subfolders.

### 2.9.1 Projects folder

To organize the projects, in the directory *COAWST/Projects/Sandy* are all the files used to simulate the Sandy case. The following files must be inside:

- Bound\_spec\_command
- coastline.mat
- coupling\_sandy.in
- create\_sandy\_application.m
- hycom\_info.mat
- ijcoast.mat
- multi\_1.at\_10m.dp.201210.grb2
- multi\_1.at\_10m.hs.201210.grb2
- multi\_1.at\_10m.tp.201210.grb2
- namelist.input
- ocean\_sandy.in
- roms\_master\_climatology\_sandy.m
- roms\_narr\_Oct2012.nc
- roms\_narr\_ref3\_Oct2012.nc
- Rweights.txt

- Sandy\_bdy.nc
- Sandy\_clm.nc
- Sandy\_clm\_ref3.nc
- sandy.h
- Sandy\_ini.nc
- Sandy\_ini\_ref3.nc
- Sandy\_init.hot
- Sandy\_ref3\_init.hot
- Sandy\_roms\_contact.nc
- Sandy\_roms\_grid.nc
- Sandy\_roms\_grid\_ref3.nc
- Sandy\_swan\_bathy.bot
- Sandy\_swan\_bathy\_ref3.bot
- Sandy\_swan\_coord.grd
- Sandy\_swan\_coord\_ref3.grd
- scrip\_sandy\_moving.nc
- scrip\_sandy\_static.nc
- specpts.mat
- swan\_narr\_Oct2012.dat
- swan\_narr\_ref3\_Oct2012.dat
- swan\_sandy.in
- swan\_sandy\_ref3.in
- tide\_forc\_Sandy.nc
- TPAR10.txt
- TPAR11.txt
- TPAR12.txt
- TPAR13.txt
- TPAR14.txt
- TPAR15.txt
- TPAR16.txt
- TPAR17.txt
- TPAR18.txt
- TPAR1.txt
- TPAR2.txt
- TPAR3.txt
- TPAR4.txt
- TPAR5.txt
- TPAR6.txt
- TPAR7.txt
- TPAR8.txt
- TPAR9.txt
- USeast\_bathy.mat
- Uweights.txt
- Vweights.txt
- wrfbdy\_d01
- wrfinput\_d01
- wrfinput\_d02
- wrflowinp\_d01

- wrflowinp\_d02

### 2.9.2 Work folder

To make the management of simulations easier, it is suggested, for each new user, the creation of the *Work* folder in main COAWST directory, with each project inserted separately within it. It is in this folder that will be simulated the test case.

```

1 cd /scratch/name.surname/COAWST
2 mkdir Work
3 cd Work
4 mkdir Sandy

```

The folder */scratch/name.surname/COAWST/Work/Sandy* must contain the following files.

- run\_sandy.sh
- limpa.sh
- link.sh

The file *run\_sandy.sh* will be used to start the simulation, *link.sh* creates symbolic links in the folder *Work*, that will be used by the models, and *limpa.sh* is used to clear the folder if an error occurs and a new integration needs to be started. The files are inside the folder */repositorio/work\_coawst*.

Therefore:

```

1 cd /scratch/name.surname/repositorio/work_coawst
2 cp limpa.sh link.sh run_sandy.sh /scratch/name.surname/COAWST/Work/Sandy

```

Go to the directory */scratch/name.surname/COAWST/Work/Sandy* and open the file *run\_sandy.sh*:

```

1 cd /scratch/name.surname/COAWST/Work/Sandy
2 nedit run_sandy.sh

```

Search for the following commands and modify the directories according to your username:

```

1 PBS -o /scratch/name.surname/COAWST/Work/Sandy/rws_total.out
2 ROOTDIR=/scratch/name.surname/COAWST

```

### 2.9.3 Compiling the test case

Go to the Sandy project folder:

```
1 /home/name.surname/COAWST/Projects/Sandy
```

Open the following files to make changes to the next steps:

- coupling\_sandy.in
- swan\_sandy.in
- swan\_sandy\_ref3.in
- ocean\_sandy.in

```
1 nedit coupling_sandy.in swan_sandy.in swan_sandy_ref3.in ocean_sandy.in
```

Look for the command line below in the file *coupling\_sandy.in*:

```
1 OCN_name = Projects/Sandy/ocean_sandy.in
```

And replace with:

```
1 OCN_name = /scratch/name.surname/COAWST/Projects/Sandy/ocean_sandy.in
```

In the files *swan\_sandy.in* and *swan\_sandy\_ref3.in* complete all directory paths below:

Modify:

```
1 READGRID COORDINATES 1 'Projects/Sandy/Sandy_swan_coord.grd' 4 0 0 FREE
2 READINP BOTTOM 1 'Projects/Sandy/Sandy_swan_bathy.bot' 4 0 FREE
3 &READINP WIND 1 'Projects/Sandy/swan_nammarr_30Sep10Nov2012.dat' 4 0 FREE
4 INITIAL HOTSTART SINGLE 'Projects/Sandy/Sandy_init.hot'
```

By:

```
1 READGRID COORDINATES 1 '/scratch/name.surname/COAWST/Projects/Sandy/Sandy_swan_coord.grd' 4 0 0 FREE
2 READINP BOTTOM 1 '/scratch/name.surname/COAWST/Projects/Sandy/Sandy_swan_bathy.bot' 4 0 FREE
3 &READINP WIND 1 '/scratch/name.surname/COAWST/Projects/Sandy/swan_nammarr_30Sep10Nov2012.dat' 4 0 FREE
4 INITIAL HOTSTART SINGLE '/scratch/name.surname/COAWST/Projects/Sandy/Sandy_init.hot'
```

In the file *ocean\_sandy.in* look for the command lines like below:

```

1 MyAppCPP = SANDY
2 VARNAME  = ROMS/External/varinfo.dat
3 GRDNAME == Projects/Sandy/Sandy_roms_grid.nc \
             Projects/Sandy/Sandy_roms_grid_ref3.nc
4 ININAME == Projects/Sandy/Sandy_ini.nc \
             Projects/Sandy/Sandy_ini_ref3.nc
5 NGCNAME =  Projects/Sandy/Sandy_roms_contact.nc
6 BRYNAME == Projects/Sandy/Sandy_bdy.nc
7 FRCNAME == Projects/Sandy/roms_narr_Oct2012.nc \
             Projects/Sandy/roms_narr_ref3_Oct2012.nc
10

```

And replace with:

```

1 MyAppCPP = Sandy
2 VARNAME  = /scratch/name.surname/COAWST/ROMS/External/varinfo.dat
3 GRDNAME == /scratch/name.surname/COAWST/Projects/Sandy/Sandy_roms_grid.nc \
             /scratch/name.surname/COAWST/Projects/Sandy/Sandy_roms_grid_ref3.nc
4 ININAME == /scratch/name.surname/COAWST/Projects/Sandy/Sandy_ini.nc \
             /scratch/name.surname/COAWST/Projects/Sandy/Sandy_ini_ref3.nc
5 NGCNAME =  /scratch/name.surname/COAWST/Projects/Sandy/Sandy_roms_contact.nc
6 BRYNAME == /scratch/name.surname/COAWST/Projects/Sandy/Sandy_bdy.nc
7 FRCNAME == /scratch/name.surname/COAWST/Projects/Sandy/roms_narr_Oct2012.nc \
             /scratch/name.surname/COAWST/Projects/Sandy/roms_narr_ref3_Oct2012.nc
10

```

Go back to the main COAWST folder and open the file *coawst.bash*:

```

1 cd /scratch/name.surname/COAWST
2 nedit coawst.bash

```

Search for the following commands and modify, if necessary:

```

1 export COAWST_APPLICATION=JOE_TC
2 export MY_ROOT_DIR=${HOME}/COAWST
3 export MY_HEADER_DIR=${MY_PROJECT_DIR}/Projects/JOE_TC

```

By:

```

1 export COAWST_APPLICATION=Sandy
2 export MY_ROOT_DIR=${HOME}/COAWST
3 export MY_HEADER_DIR=${MY_PROJECT_DIR}/Projects/Sandy

```

Activate the modules again into the file *setup\_pgi.sh*, and then compile the project with the command:

```
1 ./coawst.bash -j 4 1> coawst.pgi.sandy 2>&1 &
```

This command will create the text file *coawst.pgi.sandy* where you can follow the progress of the compilation. Open the text file with the following command and look for the final message, as shown in Figure 2.4.

```
1 nedit coawst.pgi.sandy
```

```
IPA: no IPA optimizations for 5 source files
IPA: Recompiling ./Build/get_sparse_matrix.o: new IPA information
IPA: Recompiling ./Build/master.o: new IPA information
IPA: Recompiling ./Build/mct_coupler_utils.o: new IPA information
IPA: Recompiling ./Build/mod_coupler_iounits.o: new IPA information
IPA: Recompiling ./Build/ocean_control.o: new IPA information
IPA: Recompiling ./Build/ocean_coupler.o: new IPA information
IPA: Recompiling ./Build/read_coawst_par.o: new IPA information
IPA: Recompiling ./Build/read_model_inputs.o: new IPA information
IPA: Recompiling ./Build/roms_export.o: new IPA information
IPA: Recompiling ./Build/roms_import.o: new IPA information
```

**Figure 2.4.** Final message after compiling COAWST.

If you want to follow the progress of the compilation through the terminal, use the command:

```
1 tail -f coawst.pgi.sandy
```

**WARNING** The compilation process is long!

We finished! In the main COAWST directory, a file called *coawstM* will be created. In this file all the information about your project will be compiled. Now with COAWST compiled, we will start the test case.

## 2.10 Simulating the Sandy Test Case

To simulate the case, search for the *run\_sandy.sh* file in *Work/Sandy*. Type it:

```
1 cd /scratch/name.surname/COAWST/Work/Sandy
2 nedit run_sandy.sh
```

**WARNING** The *Work* folder should contain the files *clean.sh*, *link.sh* and *run\_sandy.sh*. They can be found in the folder used as a repository. See Section 2.9.2.

When opening the file, check if the directories are in accordance with your username and type the command below to start the integration.

```
1 qsub run_sandy.sh
```

**WARNING**

The command *qsub* will submit you job. It will send the executed script to a batch of the cluster, reserving a part of the processors for your simulation.

The process will generate two files to follow the evolution of the simulation: *log.out* and *log.err*. To open, use:

```
1 nedit log.out log.err
```

Or look directly at the terminal with the command:

```
1 tail -f log.out
```

The outputs of the simulations will be stored in the *Work/Sandy* folder. If an error occurs, clean the desktop with the command:

```
1 ./limpa.sh
```

**WARNING**

The simulation of the may take several hours when integrating using 3 processors.



## 3. COAWST features

So far we have learned how to use the Kerana cluster and how to compile and run a simple COAWST test case. From now on we will enter into the specifics of the models, such as changing the number of processors and the rate of exchange of information between them.

### 3.1 Structural model files

As you may have noticed when simulating the Sandy test case, the models have files that assist the user about how to setup the project

ROMS uses the *sandy.h* file as a file containing the C pre-processing options that define the project. Access the website <https://www.myroms.org/wiki/cppdefs.h> to know more about the available options into this file. ROMS also uses the *ocean\_sandy.in* as standard input for running the model. This file defines the spatial dimensions of the project and parameters that are not informed during compilation, such as time steps, coefficients and physical constants, configuration of vertical coordinates, flags for control the output frequency, among other factors. You can learn more about this file by accessing the site <https://www.myroms.org/wiki/ocean.in>.

WRF uses the file *namelist.input* to manage information about the project, as well as physical parameterization schemes that will be used. To learn about the file description, visit [https://esrl.noaa.gov/gsd/wrfportal/namelist\\_input\\_options.html](https://esrl.noaa.gov/gsd/wrfportal/namelist_input_options.html). To learn about the physical options of the model and the references of each one, visit [http://www2.mmm.ucar.edu/wrf/users/phys\\_references.html](http://www2.mmm.ucar.edu/wrf/users/phys_references.html).

SWAN uses the file *swan\_sandy.in* as a manager. It describes several parameters, such as the project description, the input data, the grid and the initial and boundary conditions, the wave physics parameterizations, among others. To learn more about configuring the file, visit the *User Manual* tab in the website <http://swanmodel.sourceforge.net/> and look for the section *Description of commands*.

## 3.2 Modifying the number of processors

### 3.2.1 ROMS

The processors are located in the *ocean.in* file, which is located in the *Projects* folder, and are called *NtileI* and *NtileJ*. An example is in Figure 3.1. In this case, ROMS will reserve 320 processors to run, since  $16 \times 20 = 320$ .

```
! Domain decomposition parameters for serial, distributed-memory or
! shared-memory configurations used to determine tile horizontal range
! indices (Istr,Iend) and (Jstr,Jend), [1:Ngrids].
NtileI == 16                                ! I-direction partition
NtileJ == 20                                ! J-direction partition
```

**Figure 3.1.** Representation of the number of processors used in ROMS.

### 3.2.2 WRF

To change the number of WRF processors, it is necessary to modify the file *namelist.input*. You will find the variables *nproc\_x* and *nproc\_y*, as in Figure 3.2. In this case, 320 processors will be reserved for the atmospheric model.

```
nproc_x = 16,
nproc_y = 20,
```

**Figure 3.2.** Representation of the number of processors for the WRF.

### 3.2.3 SWAN

The SWAN *.in* file does not discriminate the numbers of processors used, so just change the number number of processors in *coupling.in*, as will be shown in the subsection below.

### 3.2.4 COAWST

Now that the number of processors have been modified, the coupler needs to be informed about the change. This information is communicated through the file *coupling.in*. It should contain the total number of processors to be used by the ROMS, WRF and SWAN models, as in the figure 3.3:

```
! Number of parallel nodes assigned to each model in the coupled system.
! Their sum must be equal to the total number of processors.

NnodesATM = 320                                ! atmospheric model
NnodesWAV = 64                                  ! wave model
NnodesOCN = 320                                ! ocean model
```

**Figure 3.3.** Representation of the number of processors for each COAWST module.

The *NnodesATM* refers to the total number of processors used by WRF, *NnodesWAV* is the total used by SWAN and *NnodesOCN* is total used by ROMS. Just change according to the total number of processors used in the previous steps.

Finally, it is necessary to modify the total number of processors in *run.sh*, used to submit the experiment. Add the total number of processors used by the three models, following the equation 3.1:

$$TotalProc = NnodesATM + NnodesWAV + NnodesOCN \quad (3.1)$$

Where *TotalProc* is the sum of all processors used in the models.

Now open the file *run.sh*, located inside the folder *Work*, and look for the following line:

```
1 PBS -l mppwidth= 3
```

And for the line:

```
1 aprun -n 36 coawstM ./coupling.in 1> log.out 2> log.err
```

Modify the number 3 by the total number of processors used.

### 3.3 Modifying the coupling time interval between models

To modify the information exchanged between models, open the file *coupling.in* and modify the variables *TI\_ATM2WAV*, *TI\_ATM2OCN*, *TI\_WAV2ATM*, *TI\_WAV2OCN*, *TI\_OCN2WAV*, *TI\_OCN2ATM*, as in Figure 3.4.

**WARNING** The information exchange rate is defined in seconds.

```
! Time interval (seconds) between coupling of models.
TI_ATM2WAV = 900.0d0          | atmosphere to wave coupling interval
TI_ATM2OCN = 900.0d0          | atmosphere to ocean coupling interval
TI_WAV2ATM = 900.0d0          | wave to atmosphere coupling interval
TI_WAV2OCN = 900.0d0          | wave to ocean coupling interval
TI_OCN2WAV = 900.0d0          | ocean to wave coupling interval
TI_OCN2ATM = 900.0d0          | ocean to atmosphere coupling interval
```

**Figure 3.4.** Information exchange interval between models used in COAWST. In this example, the exchange will occur every 900 seconds.





## 4. Building the WRF

Now that we have learned how to simulate a test case and change the coupling rate and number of processors, we will begin the process of creating a specific grid and the initial and boundary conditions for the WRF using the NCEP Climate Forecast System Reanalysis (CFSR; [Saha et al., 2010](#)).

### 4.1 Compiling WRF in Kerana cluster

It is recommended to copy the WRF folder inside COAWST to your work root area, in order to avoid conflicts if you choose to use the WRF without coupling to the COAWST model. Therefore:

```
1 cd /scratch/name.surname/COAWST  
2 cp -r WRF /scratch/nome.dsobrenome
```

Activate the modules in the *setup\_pgi.sh* file with the command:

```
1 source setup_pgi.sh
```

Enter the copy folder of the WRF (*/home/name.surname/WRF*) and run:

```
1 ./configure
```

**WARNING** The next step can be automated. If you choose the automated compilation, see Section [2.7](#).

If you have automated the compilation process, as shown in Section 2.7, the files *real.exe*, *wrf.exe*, *tc.exe* and *ndown.exe* will be created. If you choose the manual option on the Kerana cluster, look down for the option *Cray XC CLE / Linux x86 \_64, Cray compiler with gcc (dmpar)*.

In the next option, the following message will appear:

```
1 Compile for nesting? (1=basic, 2=preset moves, 3=vortex following) [default 1]:
```

Choose option 1.

**WARNING** End of the automated build process.

Type:

```
1 compile em_real
```

The *real.exe*, *wrf.exe*, *tc.exe* and *ndown.exe* will be generated in the */home/name.surname/WRF/test/em\_real*.

## 4.2 WRF Preprocessing System (WPS)

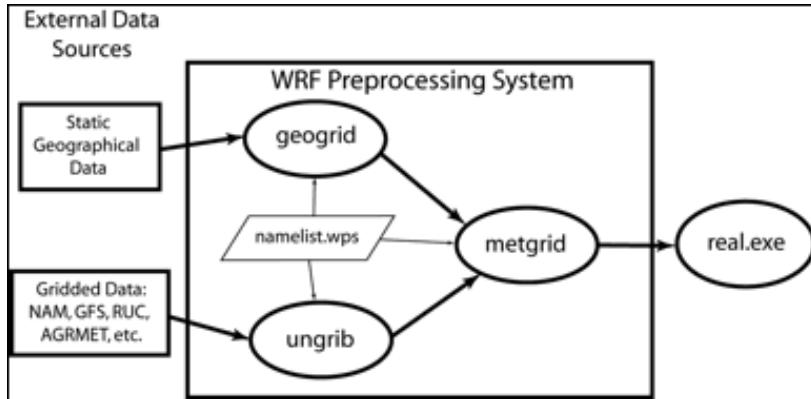
To build the initial and boundary conditions of the WRF, we will use WRF Preprocessing System (WPS). The WPS is a set of three programs that prepare the input data for the simulations. It consists of the following modules:

- ***geogrid***: Defines the model domain and interpolates the geographic data for the grid;
- ***ungrib***: Extract weather fields from *.grib* files;
- ***metgrid***: Interpolates the weather fields extracted by ungrid to the model grid defined by metgrid.

WPS is controlled by the *namelist.wps* file and is outlined in Figure 4.1.

For more information about WPS, visit: <http://www2.mmm.ucar.edu/wrf/OnLineTutorial/>.

After creating these files, the *real.exe* is used to interpolate the meteorological fields to  $\eta$ level.

**Figure 4.1.** WPS schematic.

Author: Duda (2008).

### 4.2.1 Downloading WPS

WPS is included in the COAWST package, however WPS can be downloaded by visiting [http://www2.mmm.ucar.edu/wrf/users/download/get\\_source.html](http://www2.mmm.ucar.edu/wrf/users/download/get_source.html). Download it and add the *tar.gz* file inside the cluster with *sftp*:

```

1 sftp -P2000 name.surname@acesso-hpc.cptec.inpe.br
2 cd COAWST
3 put WPSV3.9.0.1.tar.gz
  
```

To unzip, use *ssh* to enter the cluster and type:

```

1 ssh -Y name.surname@acesso-hpc.cptec.inpe.br -p 2000
2 cd COAWST
3 tar -xvzf WPSV3.9.0.1.tar.gz
  
```

### 4.2.2 Compiling WPS in the Kerana cluster

**WARNING**

As stated earlier in Section 4.2.2, WPS is included with COAWST. We recommend to use this version.

To compile the WPS and generate the executables, it is necessary to activate the libraries with *setup\_pgi.sh*, found in the directory */home/name.surname/repository*. Activate them with the command:

```

1 source setup_pgi.sh
  
```

Inside the WPS folder, enter the following command:

```
1 ./configure
```

You will be asked which machine and compiler will be used. Choose the option *Cray XE / XC CLE / Linux x86 \_64, PGI compiler (serial)*. In this case, a message will be generated that Fortran is not compatible with C and NetCDF, but ignore it.

Open *configure.wps* file:

```
1 nedit configure.wps
```

Modify:

```
1 WRF_DIR  = ../WRFV3
2 SFC       = ftn
3 SCC       = gcc
4 DM_CC    = cc
5 DM_FC    = ftn
6 FFLAGS   = -N255 -f free -h byteswapi
7 F77FLAGS = -N255 -f fixed -h byteswapi
```

By:

```
1 WRF_DIR  = /scratch/name.surname/WRF
2 SFC       = ftn
3 SCC       = gcc
4 DM_CC    = gcc
5 DM_FC    = ftn
6 FFLAGS   =
7 F77FLAGS =
```

Save the changes and start the compilation with the command:

```
1 ./compile
```

You should generate the executables *metgrid.exe* and *geogrid.exe*. However, it is possible that *ungrib.exe* is not generated. In this case, repeat:

```
1 ./configure
```

Choose the option *Linux x86\_64, PGI compiler (dmpar)*.

Open *configure.wps* file again, modify:

```
1 WRF_DIR = ../WRFV3
2 SFC      = pgf90
3 SCC      = pgcc
4 DMFC     = mpif90
5 DMCC     = mpicc
```

By:

```
1 WRF_DIR = /home/name.surname/WRF
2 SFC      = ftn
3 SCC      = gcc
4 DMFC     = ftn
5 DMCC     = gcc
```

The installer will generate only the *ungrb.exe*.

### 4.3 Building your project using CFSR

A database as option to use for simulations is the NCEP Climate Forecast System Reanalysis (CFSR) (<http://rda.ucar.edu/datasets/ds093.0>). This database is a reanalysis, from January 1979 to December 2010, where a coupled model (atmosphere-ocean-land surface and sea ice) assimilates orbital data, research cruises, oceanographic buoys and weather stations. This database has vertical resolution with 38 levels, extending from the surface to 1 hPa, with a 6-hour temporal resolution and horizontal of 0.5° for pressure levels data and 0.312° for variables on the surface.

To access the data it is necessary to register on the site. After doing this, enter the database link, click on the *Data Access* and then, *Get a subset*.

On the next page, as seen in Figure 4.2, write the temporal selection, and select in *Parameter presets* one of the three parameters of the WRF presets: (*Pressure*, *Surface* and *SST*).

**WARNING**

Only one parameter (*Pressure*, *Surface* and *SST*) can be downloaded at a time, therefore you need to repeat this step three times.

**Figure 4.2.** Screenshot of the CFSR data page.

Download the data (as *GRIB*) and separate it into three folders according to the parameters of each one: SST, Surface and Pressure. Compress them and put them in the cluster:

```

1 tar -cvzf SST.tar.gz SST/
2 tar -cvzf Pressure.tar.gz Pressure/
3 tar -cvzf Surface.tar.gz Surface/
4 ssh -Y name.surname@acesso-hpc.cptec.inpe.br -p 2000
5 mkdir Data_CFSR
6 exit
7 sftp -P2000 name.surname@acesso-hpc.cptec.inpe.br
8 cd Data_CFSR
9 put SST.tar.gz
10 put Pressure.tar.gz
11 put Surface.tar.gz
12 exit
13 ssh -Y name.surname@acesso-hpc.cptec.inpe.br -p 2000
14 cd Data_CFSR
15 tar -xvzf SST.tar.gz
16 tar -xvzf Pressure.tar.gz
17 tar -xvzf Surface.tar.gz

```

### 4.3.1 geogrid

As previously mentioned, textit geogrid generates the grid domain using geographic data. The data can be obtained from this website: [http://www2.mmm.ucar.edu/wrf/users/download/get\\_sources\\_wps\\_geog.html](http://www2.mmm.ucar.edu/wrf/users/download/get_sources_wps_geog.html), or directly from the guide repository:

```
1 /scratch/name.surname/repositorio/COAWST/WPS/geog
```

With the CFSR and *geogrid* data in hand, enter into the WPS directory and open *namelist.wps*. The file structure is shown in Figure 4.3:

```
1 nedit namelist.wps
```

```

namelist.wps - /scratch/adriano.util/COAWST3/WPS/ (em login1)
File Edit Search Preferences Shell Macro Windows
Help
namelist.wps
1 !share
2 wrf_core = 'ARW',
3 num_metgrid = 1,
4 start_date = '2004-03-15_00:00:00',
5 end_date = '2004-04-05_18:00:00',
6 interval_seconds = 21600,
7 io_form_geogrid = 2,
8 /
9
10 &geogrid
11 parent_grid_id = 1,
12 parent_grid_ratio = 1,
13 i_parent_start = 1,
14 j_parent_start = 1,
15 e_we = 600,
16 e_sn = 600,
17 geog_data_res = '10m',
18 dx = 6000,
19 dy = 6000,
20 map_proj = 'lambert',
21 ref_lat = -35,
22 ref_lon = -45,
23 trueLat1 = -35,
24 trueLat2 = -35,
25 stand_lon = -45,
26 geog_data_path = '/scratch/joao.hackerott/wrf3.7/geog',
27 opt_geogrid_tbl_path= '/scratch/adriano.util/COAWST3/WPS/geogrid'
28 /
29
30 &ungrib
31 out_formal = 'WPS',
32 prefix = 'SFC',
33 /
34
35 &netgrid
36 fg_name = 'SST','PRES','SFC',
37 io_form_metgrid = 2,
38 /

```

**Figure 4.3.** *namelist.wps* example.

Change the *geog\_data\_path* path to the directory where geographic data is located (*geog folder*), *geog\_data\_res* if you want to use other geographic data and *opt\_geogrid\_tbl\_path*, which is the *geogrid* directory.

The WPS uses a ground zero point defined by the user as the latitude and longitude degrees in *ref\_lat* and *ref\_lon*. It will generate the grid from that point and through the chosen spatial resolution. Take as example the Figure 4.3: a simulation will be prepared with 6 km of spatial resolution (*dx* and *dy* = 6000), in Lambert projection, with 600 grid points. Using as reference latitude -35°longitude -45°, a grid will be created with 600 points in the west-east direction (*e\_we*) and 600 points in the south-north direction (*e\_sn*), with each point having a spacing of 6000 meters (6 km) between them.

To start *geogrid*, you need a file with a Shell extension (.sh), which submits program to be executed in the cluster. It can be found in the */repository/WPS\_scripts* directory. Move the contents of the folder to the WPS directory with the commands:

```
1 cd /home/name.surname/repositorio/WPS_scripts
2 mv qsub_geogrid.sh qsub_ungrib.sh qsub_metgrid.sh /home/name.surname/COAWST/WPS
```

The structure of the *qsub\_geodrid.sh* file is shown in Figure 4.4. Modify the directory according to your user.

```
#!/bin/sh
#PBS -l mppwidth=1
#PBS -N GEOGRID
#PBS -j oe
#PBS -o Joe_test.out
#PBS -l walltime=02:30:00
#PBS -q workq
echo "Running GEOGRID on KERANA"
#
export MPICH_ENV_DISPLAY=1
export MPICH_ABORT_ON_ERROR=1
export MPICH_RANK_REORDER_DISPLAY=1
export MPICH_RANK_REORDER_METHOD=1
export MALLOC_MMAP_MAX_=0
export MALLOC_TRIM_THRESHOLD_=536870912
export OMP_NUM_THREADS=1
#
EXECDIR=/scratch/nome.sobrenome/COAWST/WPS
cd $EXECDIR
aprun -n 1 geogrid/src/geogrid.exe 1> log.out 2> log.err
\rm geogrid.log.00*
```

**Figure 4.4.** *qsub\_geodrid.sh* example.

To run *geogrid*, use *qsub* command:

```
1 qsub qsub_geodrid.sh
```

Two files will be generated: *log.err* e *log.out*. you can follow how the program is being executed and search for errors. The program completion message is shown in Figure 4.5, in the file *log.out*.

```
!!!!!!!!!!!!!!!
! Successful completion of geogrid.
!!!!!!!!!!!!!!!
Application 593294 resources: utime ~813s, stime ~1s
```

**Figure 4.5.** Example of completion of *geogrid*.

When completing the execution of *geogrid*, the file *geo\_em\_d01.nc* will be generated. You can view the NetCDF file with Ncview:

```
1 module load netcdf
2 ncview geo_em_d01.nc
```

You will always need to run *geogrid* until you find a domain that is suitable for your project, since WPS does not have a graphical interface. One solution to work around the problem is to use NCAR Command Language version 6.4 (NCL; UCAR, 2017) to plot the image of the chosen domain. This will be the subject of the next subsection.

### 4.3.2 Using the NCL to plot the grid domain

To install the NCL, move the *NCL* folder inside the repository:

```
1 mv /scratch/name.surname/repositorio/NCL /scratch/name.surname
```

Open the *.bashrc* file:

```
1 cd
2 nedit .bashrc
```

Add the following command lines, changing to your username:

```
1 export NCARG_ROOT=/scratch/name.surname/NCL
2 export PATH=$NCARG_ROOT/bin:$PATH
```

Save the file, reload the modules inside *.bashrc* and then execute NCL.

```
1 source .bashrc
2 ncl
```

Check if the NCL has been initialized as in Figure 4.6.

```
Copyright (C) 1995-2015 - All Rights Reserved
University Corporation for Atmospheric Research
NCAR Command Language Version 6.3.0
The use of this software is governed by a License Agreement.
See http://www.ncl.ucar.edu/ for more details.
ncl 0>
ncl 1>
```

**Figure 4.6.** Example of NCL initialization.

Enter the repository and open the *plotgrids.ncl* file. This script generates the domain image from the data in the *namelist.wps* file.

```
1 cd /scratch/name.surname/repositorio  
2 nedit plotgrids.ncl
```

Modify the *namelist.wps* directory according to your user area:

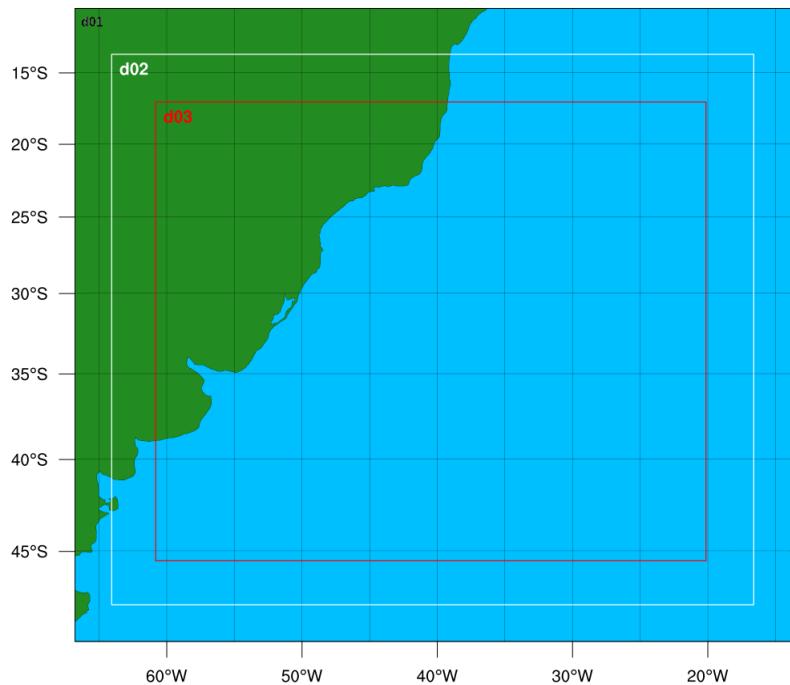
```
1 filename = "/home/name.surname/COAWST/WPS/namelist.wps"
```

Run the code with the command:

```
1 ncl plotgrids.ncl
```

The code will search for the grid points and the domain resolution through the file *namelist.wps* and show it on the screen, as shown in Figure 4.7. Repeat the process until you find a domain that suits your needs.

### WPS Domain Configuration



**Figure 4.7.** Figure generated by the *plotgrids.ncl* file. In this example, three domains were generated.

### 4.3.3 *ungrib*

With the domain prepared by *geogrid*, we can proceed to *ungrib*, which extracts the data from the *Grib* files. We will start creating the SST data. Open *namelist.wps* and change the *prefix* to *SST*, in the *& ungrib* section, as shown in Figure 4.8.

```
&ungrib
  out_format = 'WPS',
  prefix = 'SST',
/
```

**Figure 4.8.** *&ungrib* example.

Import the Variable Table for SST through a symbolic link. Type in the terminal:

```
1 ln -sf ungrib/Variable_Tables/Vtable.SST Vtable
```

**WARNING**

The *Vtable* is used to read the files in *Grib* format. It is mandatory to use the corresponding *Vtable* to the chosen database. You can consult them at *WPS/ungrib/Variable\_Tables*.

Create the symbolic links with the SST data through the *link\_grib.csh* file:

```
1 ./link_grib.csh /scratch/name.surname/Dados_CFSR/SST/*
```

Open the *qsub\_ungrib.sh* file and modify the path according to your user and type:

```
1 qsub qsub_ungrib.sh
```

**WARNING**

The *qsub\_ungrib.sh* file is located in the */repository/WPS\_scripts* folder.

Several files will be created with the initial *SST*: followed by the date chosen for simulation. To check if everything went well, look for the message in Figure 4.9 at the end of the file *log.out*.

```
!!!!!!Successful completion of ungrib!!!!!!
Application 593300 resources: utime ~18s, stime ~1s
```

**Figure 4.9.** Final message in the *log.out* file when running *ungrib*.

We will proceed to the CFSR surface data. Change the *prefix* in *namelist.wps* (Figure 4.8). Modify *SST* by *SFC* and save the modification.

Import the CFSR *Variable Table* with the symbolic link creation command:

```
1 ln -sf ungrib/Variable_Tables/Vtable.CFSR Vtable
```

Create the symbolic links to the CFSR surface data with the command below and run *ungrib* again:

```
1 ./link_grib.csh /scratch/name.surname/Dados_CFSR/Surface/*
2 qsub qsub_ungrib.sh
```

Look for the final message as shown in Figure 4.9.

Change the *prefix* of *namelist.wps* (Figure 4.8), replace *SFC* with *PRES* and save the file.

Create the symbolic links to the pressure data and run *ungrib*:

```
1 ./link_grib.csh /scratch/name.surname/Dados_CFSR/Pressure/*
2 qsub qsub_ungrib.sh
```

Finally, look for the final message as identified in Figure 4.9. At the end of these steps, there will be several files with the initials *SST*, *PRES*, *SFC* followed by the dates of the chosen period.

#### 4.3.4 metgrid

To interpolate the data generated by *ungrib.exe*, open *namelist.wps* and change on the *fg\_name* tab the names used in *ungrib*. In the previous case, *PRES*, *SFC* and *SST* were used, as in the example in Figure 4.10:

```
&metgrid
  fg_name = 'PRES', 'SFC', 'SST',
  io_form_metgrid = 2,
/
```

**Figure 4.10.** *&metgrid* example.

Open and then change the *qsub\_metgrid.sh* file path and run:

```
1 qsub qsub_metgrid.sh
```

Move the files generated by *metgrid* (*met\_em.d01\**) to the real WRF case directory:

```
1 mv met_em.d01* /home/name.surname/WRF/test/em_real
```

**WARNING**

If you are an experienced user, you can create symbolic links pointing to the *met* files instead of moving them.

#### 4.3.5 *real*

We will use the *real.exe* program to generate the files that will be used in WRF. Open the file *namelist.input*, which is in the WRF folder (*/home/name.surname/WRF/test/em\_real*), and modify it according to your chosen domain and simulation time in *namelist.wps*.

For the WRF to simulate the SST, it is necessary to include the following variables at the end of the section *&time\_control* (Figure 4.11):

```
1 io_form_auxinput4 = 2,  
2 auxinput4_inname = "wrfflowinp_d<domain>",  
3 auxinput4_interval = 360,
```

The *io\_form\_auxinput4* refers to the final format of the *wrfflowinp* file that will be generated by the *real*. The *auxinput4\_inname* is the name of the boundary condition file for SST and *auxinput4\_interval* is the time interval, in minutes, of the boundary file.

```

&time_control
run_days                                = 5,
run_hours                                 = 0,
run_minutes                               = 0,
run_seconds                               = 0,
start_year                                = 2008,
start_month                               = 11,
start_day                                 = 21,
start_hour                                 = 00,
start_minute                               = 00,
start_second                               = 00,
end_year                                  = 2008,
end_month                                 = 11,
end_day                                   = 25,
end_hour                                   = 18,
end_minute                                 = 00,
end_second                                 = 00,
interval_seconds                          = 21600,
input_from_file                           = .true.,
history_interval                         = 360.,
frames_per_outfile                        = 80,
restart                                    = .false.,
restart_interval                          = 50000,
io_form_history                           = 2
io_form_restart                           = 2
io_form_input                             = 2
io_form_boundary                          = 2
debug_level                               = 0
io_form_auxinput4                         = 2
auxinput4_inname                           = "wrflowinp_d<domain>"
auxinput4_interval                         = 360,
/

```

**Figure 4.11.** *&time\_control* example.

Also, in the *&physics* (Figure 4.12) in the *namelist.input* the option to update the SST:

```
| sst_update = 1,
```

```

&physics
sst_update                                = 1,
mp_physics                                 = 6,
ra_lw_physics                             = 4,
ra_sw_physics                             = 3
swint_opt                                  = 1,
sf_sfclay_physics                         = 1,
sf_surface_physics                        = 2,
bl_pbl_physics                            = 1,
cu_physics                                 = 1,
isfflx                                     = 1,
ifsnnow                                     = 1,
icloud                                      = 1,
surface_input_source                      = 1,
num_soil_layers                           = 4,
sf_urban_physics                          = 0,
sst_skin                                    = 0,

```

**Figure 4.12.** *&physics* example.

After completing the modifications, you must use the file *qsub\_real.sh* to submit the job. The file is located in

the */repository/WRF\_scripts* directory. Move the file to the *em\_real* folder, change the directory according to your user and execute:

```
1 cd /home/name.surname/repositorio/WRF_scripts
2 mv qsub_real.sh /home/name.surname/WRF/test/em_real
3 qsub qsub_real.sh
```

**WARNING** The *qsub\_real.sh* file is located in the */repository/WRF\_scripts* folder.

To check if *real* succeeded, look at the end of the *log.out* file for the success message as shown in Figure 4.13.

```
d01 2007-02-18_00:00:00 real_em: SUCCESS COMPLETE REAL_EM INIT
d01 2007-02-18_00:00:00 real_em: SUCCESS COMPLETE REAL_EM INIT
d01 2007-02-18_00:00:00 Timing for loop # 41 = 2 s.
d01 2007-02-18_00:00:00 real_em: SUCCESS COMPLETE REAL_EM INIT
d01 2007-02-18_00:00:00 real_em: SUCCESS COMPLETE REAL_EM INIT
d01 2007-02-18_00:00:00 Timing for loop # 41 = 2 s.
d01 2007-02-18_00:00:00 real_em: SUCCESS COMPLETE REAL_EM INIT
Application 593290 resources: utime ~11862s, stime ~117s
```

**Figure 4.13.** Successful completion of *real*.

Upon completing *real*, three files will be generated: *wrfbdy\_d01*, *wrfinput\_d01* and *wrflowinp\_d01*, if you have chosen only one domain. After that, the WRF conditions are ready to be copied to your project folder in COAWST:

```
1 cp wrfbdy_d01 wrfinput_d01 wrflowinp_d01 /home/name.surname/COAWST/Projects/project_name
```

## 4.4 Using the WRF

With the files ready, it is possible to start the simulation of your project using only the WRF.

To run the simulation on the cluster, it is necessary to use the file *qsub\_wrf.sh*. It is located in the */repository/WRF\_scripts* directory. Move the file to the WRF *in\_real* folder, modify the directory according to your user and save:

```
1 cd /home/name.surname/repositorio/WRF_scripts
2 mv qsub_wrf.sh /home/name.surname/WRF/test/em_real
3 nedit qsub_wrf.sh
```

To start the simulation, type:

```
! qsub qsub_wrf.sh
```

**WARNING**

The *qsub\_wrf.sh* file is located in the */repository/WRF\_scripts* folder.

Follow the evolution of the job with the *log.out* and *log.err* files. If you want to have the information updated by the terminal, type:

```
! tail -f log.out
```



## 5. Building the ROMS

### 5.1 Installing Python libraries on the personal computer

Before generating the ROMS files, it is necessary to install the libraries and modules. In this chapter, we address how to download them, as well as the commands in the terminal to install through *apt-get*.

There are two ways to install the required libraries: manually, and through Anaconda. Anaconda is a free and open source platform for Python and R and has a package manager called Conda, which makes easy to install the libraries. In this guide, I will present both options. For convenience, I suggest installing using Conda (Section 5.1.2), but beware of version conflicts, as the Conda repositories are constantly updated.

**WARNING**

ROMS files must be builded on your own computer and no longer on Kerana cluster, as previously done with WRF.

#### 5.1.1 Installing manually

**WARNING**

Consider installing libraries with Anaconda will be faster than installing manually. See Section 5.1.2.

Open the terminal, and create a folder in your *\$HOME* called *libsmodels*. The downloaded files and libraries will be placed in it.

```
1 sudo mkdir libsmodels
```

If you do not have *gfortran* and *g ++* installed, install via *apt-get*.

```

1 sudo apt-get install
2 sudo apt-get install g++

```

Next, download OpenMPI 2.0.1 (<https://www.open-mpi.org/software/ompi/v2.0/>), extract the file into the *libsmodels* folder and type in the terminal:

```

1 export F77=gfortran
2 export FC=gfortran
3 export FC=gfortran
4 export CC=gcc
5 export CXX=g++
6 export FFLAGS="-m64 -fdefault-integer-8"
7 export FCFLAGS="-m64 -fdefault-integer-8"
8 export CFLAGS=-m64
9 export CXXFLAGS=-m64

```

Enter the OpenMPI folder and type:

```

1 ./configure --prefix=$HOME/libsmodels
2 sudo make
3 sudo make check
4 sudo make install

```

Go back to your *\$HOME* and open the hidden file *.bashrc* using *gedit* and write the following in the final lines of the file:

```

1 cd $HOME
2 gedit .bashrc
3 export PATH=$HOME/libsmodels/bin:$PATH
4 export LD_LIBRARY_PATH=$HOME/libsmodels/lib:$LD_LIBRARY_PATH
5 export PATH=$HOME/libsmodels/include:$PATH

```

Save the file and enter the following command to update the terminal and apply the changes.

```

1 source .bashrc

```

Test the installation of OpenMPI is correct, type in the terminal:

```

1 ompi_info -a grep 'Fort integer size'

```

The result should be:

```
1 Fort integer size: 8
```

If you have not installed it on your computer, you will need to install the *m4*, *make* e *perl*. To check if your computer already has these programs installed:

```
1 which m4
2 which make
3 which perl
```

If not, install from the following commands:

```
1 sudo apt-get install m4
2 sudo apt-get install make
3 sudo apt-get install perl
```

Install the *szip 2.1* (<https://www.hdfgroup.org/HDF5/release/obtain5.html>). Download, unzip it in the *\$HOME/libsmodels* folder, enter the *szip* folder and type the following:

```
1 export F77=gfortran
2 export FC=gfortran
3 export CC=gcc
4 export CXX=g++
5 export FFLAGS="-m64 -fdefault-integer-8"
6 export FCFLAGS="-m64 -fdefault-integer-8"
7 export CFLAGS=-m64
8 export CXXFLAGS=-m64
9 ./configure --prefix=$HOME/libsmodels
10 sudo make
11 sudo make check
12 sudo make install
```

Install the *zlib 1.2.8* (<http://zlib.net/>). Download, unzip it in the *\$HOME/libsmodels* folder, enter the *zlib* folder and type the following:

```

1 export F77=gfortran
2 export FC=gfortran
3 export CC=gcc
4 export CXX=g++
5 export FFLAGS="-m64 -fdefault-integer-8"
6 export FCFLAGS="-m64 -fdefault-integer-8"
7 export CFLAGS=-m64
8 export CXXFLAGS=-m64
9 ./configure --prefix=$HOME/libmodels
10 sudo make
11 sudo make check
12 sudo make install

```

Install the *Curl* 7.50.3 (<http://curl.haxx.se/download.html>). Download, unzip it in the *\$HOME/libmodels* folder, enter the *Curl* folder and type the following:

```

1 export F77=gfortran
2 export FC=gfortran
3 export CC=gcc
4 export CXX=g++
5 export FFLAGS="-m64 -fdefault-integer-8"
6 export FCFLAGS="-m64 -fdefault-integer-8"
7 export CFLAGS=-m64
8 export CXXFLAGS=-m64
9 ./configure --prefix=$HOME/libmodels/
10 sudo make
11 sudo make check
12 sudo make install

```

Install the libraries needed to use HDF5. Type it:

```
1 sudo apt-get install libgtk2.0-dev
```

Install the *HDF5* 1.8.9 (<https://www.hdfgroup.org/HDF5/release/obtain5.html>). Download, unzip it in the *\$HOME/libmodels* folder, enter the *HDF5* folder and type the following:

```

1 export FC=gfortran
2 export CC=gcc
3 export CXX=g++
4 export LDFLAGS=-L$HOME/libmodels/lib
5 export CPPFLAGS=-I$HOME/libmodels/include

```

In the same directory, type:

**WARNING** Type everything from the box below in the same command line!

```
1 ./configure --enable-fortran=yes --enable-fortran2003=yes --enable-cxx=yes  
2 --with-szlib=$HOME/libsmodeles --with-zlib=$HOME/libsmodeles --enable-hl  
3 --enable-shared --prefix=$HOME/libsmodeles
```

Complete the installation of *HDF5* with the commands below:

```
1 sudo make  
2 sudo make check  
3 sudo make install
```

The next step is to install the *NetCDF-C 4.4.1* (<https://www.unidata.ucar.edu/downloads/netcdf/index.jsp>). Download, unzip it in the *\$HOME/libsmodeles* folder, enter the *NetCDF* folder and type the following:

```
1 export FC=gfortran  
2 export CC=gcc  
3 export CXX=g++  
4 export LDFLAGS=-L$HOME/libsmodeles/lib  
5 export CPPFLAGS=-I$HOME/libsmodeles/include
```

Then type:

**WARNING** Type everything from the box below in the same command line!

```
1 ./configure --prefix=$HOME/libsmodeles --enable-netcdf4 --enable-shared --enable-dap  
2 --with-hdf5=$HOME/libsmodeles
```

Complete the installation of *NetCDF-C* with the commands below:

```
1 sudo make  
2 sudo make check  
3 sudo make install
```

Install the *NetCDF-Fortran 4.2* (<https://www.unidata.ucar.edu/downloads/netcdf/index.jsp>). Download, unzip it in the *\$HOME/libsmodeles* folder, enter the *NetCDF-Fortran* folder and type the following:

```

1 export FC=gfortran
2 export F77=gfortran
3 export CC=gcc
4 export CXX=g++
5 export LDFLAGS=-L$HOME/libmodels/lib
6 export CPPFLAGS=-I$HOME/libmodels/include
7 export LIBS='-lnetcdf -lhdf5 -lhdf5_hl -lz'
```

Type:

**WARNING** Type everything from the box below in the same command line!

```

1 ./configure --prefix=$HOME/libmodels --enable-netcdf4 --enable-shared
2 --with-hdf5=$HOME/libmodels --enable-dap
```

Finish installing the *NetCDF-Fortran*:

```

1 sudo make
2 sudo make check
3 sudo make install
```

Install the *JasPer 1.900.1* (<http://www.linuxfromscratch.org/blfs/view/svn/general/jasper.html>). Download, unzip it in the *\$HOME/libmodels* folder, enter the *Jasper* folder and type the following:

```

1 export FC=gfortran
2 export F77=gfortran
3 export CC=gcc
4 export CXX=g++
5 ./configure --prefix=$HOME/libmodels
6 sudo make
7 sudo make check
8 sudo make install
```

Instal the *libpng 1.6.24* ([http://sourceforge.net/projects/libpng/?source=typ\\_redirect](http://sourceforge.net/projects/libpng/?source=typ_redirect)). Download, unzip it in the *\$HOME/libmodels* folder, enter the *libpng* folder and type the following:

```

1 export FC=gfortran
2 export F77=gfortran
3 export CC=gcc
4 export CXX=g++
5 export LDFLAGS=-L$HOME/libmodels/lib
6 export CPPFLAGS=-I$HOME/libmodels/include
7 export LIBS='-lz'
8 ./configure --with-zlib=$HOME/libmodels --prefix=$HOME/libmodels
9 sudo make
10 sudo make check
11 sudo make install
```

Install the *Udunits 2.1.24* (<ftp://ftp.unidata.ucar.edu/pub/udunits/>). Download, unzip it in the \$HOME/libsmodels folder, enter the *Udunits* folder and type the following:

```
1 export FC=gfortran
2 export F77=gfortran
3 export CC=gcc
4 export CXX=g++
5 ./configure --prefix=$HOME/libsmodels
6 sudo make
7 sudo make check
8 sudo make install
```

Install NetCDF4 for Python. In this case, download Git first and use it to download.

```
1 sudo apt-get install git
```

Clone the *NetCDF4-Python* repository with the command:

```
1 git clone https://github.com/Unidata/netcdf4-python
```

Entre na pasta do NetCDF4-Python e modifique o arquivo *setup.cfg.template* e salve:

```
1 ncconfig      = $HOME/libsmodels/bin/nc-config
2 netCDF4_dir   = $HOME/libsmodels
3 HDF5_dir     = $HOME/libsmodels
4 szip_dir     = $HOME/libsmodels
5 jpeg_dir     = $HOME/libsmodels
6 curl_dir     = $HOME/libsmodels
```

Change the file name *setup.cfg.template* to *setup.cfg*:

```
1 mv setup.cfg.template setup.cfg
```

Create inside the *libsmodels* folder, the *lib/python2.7/site-packages* subfolders:

```
1 mkdir lib
2 cd lib
3 mkdir python2.7
4 cd python2.7
5 mkdir site-packages
```

Go back to the *NetCDF4-Python* main folder and compile:

```
1 python setup.py build
2 python setup.py install --prefix=$HOME/libsmmodels
```

If you want to check if the installation was successful, try to import the library into Python. Type in the terminal:

```
1 python
2 import netCDF4
```

Install the *mpi4py* (<https://pypi.python.org/pypi/mpi4py>). Extract and type:

```
1 python setup.py build
2 python setup.py install --prefix=$HOME/libsmmodels
```

To test the installation, type:

```
1 mpiexec -n 5 python demo/helloworld.py
```

Install the *ESMF* e *ESMPy* (<https://www.earthsystemcog.org/projects/esmpy/releases>). First, type the following commands to install the necessary library files:

```
1 sudo apt-get install python-setuptools
2 sudo easy_install pip
3 sudo pip install distribute
4 sudo pip install nose
```

Inside the ESMF folder, type in the terminal:

```
1 export ESMF_F90COMPILER=gfortran
2 export ESMF_NETCDF_LIBS="-lncdf -lnetcdff"
3 export ESMF_DIR=$HOME/libsmmodels/packages/esmp_ESMF_6_3_0rp1_ESMP_01/esmf
4 export ESMF_NETCDF="local"
5 export ESMF_COMPILER=gfortran
6 export ESMF_COMM=openmpi
7 export ESMF_TESTEXHAUSTIVE=on
8 export ESMF_TESTSHAREDOBJ=on
9 export ESMF_NETCDF_INCLUDE=$HOME/libsmmodels/include
10 export ESMF_NETCDF_LIBPATH=$HOME/libsmmodels/lib
11 export ESMF_INSTALL_PREFIX=$HOME/libsmmodels/ESMF
12 make
13 make check
14 make all_tests
15 make install
```

Enter in the *scr/addon/ESMPy* folder:

```
1 cd src/addon/ESMPy/
```

Type:

**WARNING** Type everything from the box below in the same command line!

```
1 python setup.py build  
2 --ESMFMKFILE=$HOME/libmodels/ESMF/lib/lib0/Linux.gfortran.64.openmpi.default/esmf.mk
```

Complete the installation:

```
1 python setup.py install --prefix=$HOME/libmodels  
2 python setup.py test
```

Install Octant, using Git:

```
1 git clone https://github.com/hetland/octant
```

In the *octant/external* folder, delete everything and download the following files:

```
1 rm -rf *  
2 git clone https://github.com/sakov/gridgen-c.git  
3 git clone https://github.com/sakov/gridutils-c  
4 git clone https://github.com/sakov/csa-c  
5 git clone https://github.com/sakov/nanoc
```

Rename the folders to, respectivamente: *gridgen*, *gridutils*, *csa* and *nn*. Enter the *nn/nn* folder, export the libraries and install *nn*:

```
1 cd nn/nn/  
2 export FC=gfortran  
3 export F77=gfortran  
4 export CC=gcc  
5 export CXX=g++  
6 ./configure --prefix=$HOME/libmodels  
7 make  
8 make install
```

Install *csa* in the same way as *nn* was installed:

```

1 cd ../../csa/csa/
2 export FC=gfortran
3 export F77=gfortran
4 export CC=gcc
5 export CXX=g++
6 ./configure --prefix=$HOME/libsmodels
7 make
8 make install

```

In the *csa* folder, copy the *csa.o* file to the *\$HOME/libsmodels/lib* folder:

```
1 cp csa.o $HOME/libsmodels/lib
```

Install the *gridutils*:

```

1 cd ../../gridutils/gridutils/
2 export FC=gfortran
3 export F77=gfortran
4 export CC=gcc
5 export CXX=g++
6 LDFLAGS=-L$HOME/libsmodels/lib
7 ./configure --prefix=$HOME/libsmodels CPPFLAGS=-I$HOME/libsmodels/include
8 make
9 make install

```

Install the *gridgen*:

```

1 cd ../../gridgen/
2 export FC=gfortran
3 export F77=gfortran
4 export CC=gcc
5 export CXX=g++
6 LDFLAGS=-L$HOME/libsmodels/lib
7 ./configure -prefix=$HOME/libsmodels CPPFLAGS=-I$HOME/libsmodels/include

```

Open the file *makefile* and change the file as shown below:

```

1 CFLAGS = -g -O2 -Wall -I$HOME/libsmodels/include
2 LDFLAGS = -L$HOME/libsmodels/lib
3 LIBS = -lgm -lm -L$HOME/libsmodels/lib

```

Save the file, close and type in the terminal:

```
1 make  
2 make lib  
3 make shlib  
4 make install
```

Inside the folder `$HOME/libmodels/octant/octant`, modify the file `grid.py` like the example below (it is approximately on row 898).

**WARNING** Pay close attention to the Python indentation inside the files!

```
1 self._libgridgen = np.ctypeslib.load_library('libgridgen.so', '$HOME/libmodels/lib')  
2 print octant.__path__[0]  
3 self._libgridgen = np.ctypeslib.load_library('_gridgen', octant.__path__[0])
```

Salve e feche o arquivo. Digite:

```
1 python setup.py build --fcompiler=gfortran
```

In `$HOME/libmodels/octant/octant`, open the file `csa.py` and modify as below:

```
1 _csa = np.ctypeslib.load_library('csa.o', '$HOME/libmodels/libs')
```

Finally, in `$HOME/libmodels/octant`, type:

```
1 python setup.py install --prefix=$HOME/libmodels/octant
```

Export the paths in `.bashrc`. Remember the: `.bashrc` file is in your `$HOME`.

```
1 export PYTHONPATH=$PYTHONPATH:$HOME/libmodels/octant  
2 export PYTHONPATH=$PYTHONPATH:$HOME/libmodels/octant/include  
3 export PYTHONPATH=$PYTHONPATH:$HOME/libmodels/octant/lib/python2.7/site-packages  
4 export PYTHONPATH=$PYTHONPATH:$HOME/libmodels/octant/lib
```

Update the terminal with the new `.bashrc`. Type:

```
1 source .bashrc
```

### 5.1.2 Using Conda to install ROMS libraries

Anaconda is a distribution for Python and R that supports the various libraries and packages used in this guide, making installation easier. To download Anaconda go to <https://www.continuum.io/downloads>. To install, just enter the directory where the installation file is located, change the installation permissions of the file and follow the installer's instructions:

```
1 sudo chmod 770 Anaconda2.sh  
2 ./Anaconda2.sh
```

Next, enter the following commands on the terminal to install the libraries needed to build a project on ROMS using Python:

```
1 conda install -c conda-forge openmpi  
2 conda install -c anaconda gcc  
3 conda install -c mutirri szip  
4 conda install -c anaconda zlib  
5 conda install -c anaconda curl  
6 conda install -c anaconda hdf5  
7 conda install -c conda-forge netcdf-fortran  
8 conda install -c scitools jasper  
9 conda install -c anaconda libpng  
10 conda install -c conda-forge udunits  
11 conda install -c anaconda netcdf4  
12 conda install -c anaconda cython  
13 conda install -c anaconda numpy  
14 conda install -c anaconda scipy  
15 conda install -c anaconda mpi4py  
16 conda install -c conda-forge esmf
```

## 5.2 Installing Pyroms

With the libraries already installed, this section will show you how to install Pyroms. It is a collection of tools to create ROMS files. It was originally started by Rob Hetland as a project on GoogleCode and rewritten by Frederic Castruccio.

Download PyROMS inside your *libsmodels* folder:

```
1 git clone https://github.com/ESMG/pyroms.git
```

Enter the directory *\$HOME/libsmodels/pyroms/pyroms\_toolbox* and type.

```
1 export FC=gfortran
2 export F77=gfortran
3 export CC=gcc
4 export CXX=g++
5 python setup.py build --fcompiler=gfortran
6 python setup.py install --prefix=$HOME/libsmmodels/pyroms
```

Install the bathy\_smoothen:

```
1 cd ../../bathy_smoothen/external/lp_solve_5.5/lp_solve/
2 sh ccc
3 cp bin/ux64/lp_solve $HOME/libsmmodels/bin/
4 cd ../../lpsolve55
5 sh ccc
6 cp bin/ux64/* $HOME/libsmmodels/lib/
```

Enter the following directory and open the *setup.py*:

```
1 cd ../../lp_solve_5.5/extra/Python/
2 gedit setup.py
```

Change the paths as described below, in the file *setup.py*. Remember the indentation!

```
1 include_dirs=['..','/usr/include','$HOME/libsmmodels/include'],
2 library_dirs=['$HOME/libsmmodels/pyroms/lib/','$HOME/libsmmodels/lib'],
```

Install:

```
1 python setup.py build
2 python setup.py install --prefix=$HOME/libsmmodels/pyroms
```

Enter the bathy\_smoothen directory and type in the terminal:

```
1 cd ../../../../../../bathy_smoothen/
2 export FC=gfortran
3 export F77=gfortran
4 export CC=gcc
5 export CXX=g++
6 python setup.py build --fcompiler=gfortran
7 python setup.py install --prefix=$HOME/libsmmodels/pyroms
```

Install Pyroms.

```
1 cd ../pyroms/pyroms/
2 python setup.py install --prefix=$HOME/libsmodels/pyroms
```

Open the file *hgrid.py* and change the line 1028 as below:

```
1 self._libgridgen = np.ctypeslib.load_library('libgridgen.so', '$HOME/libsmodels/lib')
```

In the file *hgrid.py*, add the following library at line 21:

```
1 from numpy import amin
```

Save, close *hgrid.py* and type:

```
1 cd ../
2 export FC=gfortran
3 export F77=gfortran
4 export CC=gcc
5 export CXX=g++
6 python setup.py build --fcompiler=gfortran
7 python setup.py install --prefix=$HOME/libsmodels/pyroms
```

Coloque os caminhos no seu *.bashrc*:

```
1 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/lib/python2.7/site-packages
2 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/lib/python2.7/site-packages
3 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/lib/python2.7/site-packages
4 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/lib
5 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/lib/python2.7/site-packages/pyroms_toolbox/Grid_HYCOM
6 export PYTHONPATH=$PYTHONPATH:$HOME/libsmodels/pyroms/examples
```

Update the terminal with the new *.bashrc*:

```
1 source .bashrc
```

Go to *\$HOME/libsmodels/pyroms/lib/python2.7/site-packages/bathy\_smoother/* and delete the file *lpsolve55.so*.

Go to the *pyroms\_toolbox/* folder and open *\_\_init\_\_.py* file. Add a comment or erase the content on the row that imports the *from move\_river\_t import move\_river\_t* (row 51). Then, in the folder *Grid\_HYCOM*, open the file *flood\_fast.py* and comment or erase the content on row 9 (*import creep*).

Save and import pyroms\_toolbox using Python:

```
1 python
2 import pyroms_toolbox
```

### 5.3 The model2roms toolbox

In this section, you will be introduced to install model2roms, the toolbox used to generate ROMS forcing files and its grid. It was created and is being updated by Trond Kristiansen ([Kristiansen, 2019](#)) and we adapted to the needs of LOA-INPE.

To download *model2roms*, enter *home* on your computer and download using Github:

```
1 cd $HOME
2 git clone https://github.com/usutil/model2roms
```

Change the folder name from *model2roms-master* to *model2roms*:

```
1 mv model2roms-master model2roms
```

#### 5.3.1 Introduction to the ROMS grid

To generate the ROMS grid, we will use the code *make\_roms\_grid.py*, located in the *model2roms/grid* directory. The code supports the SRTM30\_plus data ([Becker et al., 2009](#)), however, for practical purposes, we will exemplify only the data from ETOPO1 ([Amante & Eakins, 2009](#)), which will be discussed in the next subsection.

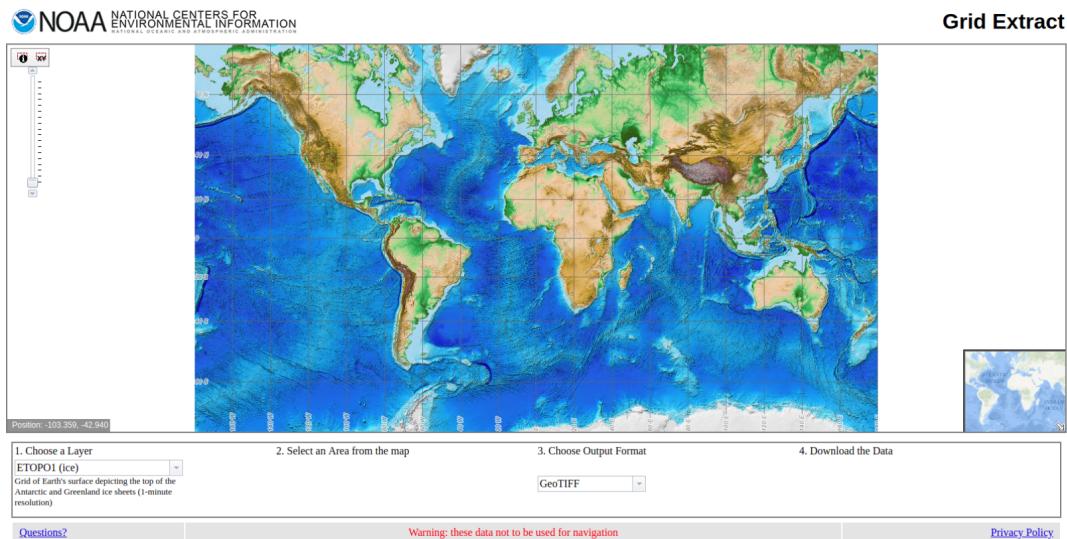
Install basemap to use the code, if you have not previously installed

```
1 conda install -c anaconda basemap
```

#### 5.3.2 ETOPO1 Arc-Minute Global Relief Model

The ETOPO1 ([Amante & Eakins, 2009](#); Figure 5.1) is produced by the National Geophysical Data Center and provides two layers of relief information. The layers include bathymetry over the oceans and some of the main lakes on Earth. Terrestrial topography and ocean bathymetry are based on SRTM30 topography and through various bathymetric cruises

Visit the site <https://www.ngdc.noaa.gov/mgg/global/global.html> and subset your area of interest on the ETOPO1 website map and download the data in NetCDF format.



**Figure 5.1.** ETOPO1 webpage.

### WARNING

The script *make\_roms\_grid.py* does not allow the selection of the latitude and longitude limits of the grid, therefore, be as accurate as possible when subsetting of your area.

### Creating the ROMS grid

Enter the *model2roms* directory and look for the *grid* subfolder. Open the code *make\_roms\_grid.py*.

```
1 cd $HOME/model2roms/grid
2 gedit make_roms_grid.py
```

As Figure 5.2 shown, the variables that will be modified are:

- **grd\_name:** Grid name;
- **grd\_final:** The NetCDF file name of the grid;
- **etopo1\_dir:** The directory where the ETOPO1 NetCDF file is located;
- **srtm\_dir:** The directory where the SRTM\_30\_PLUS NetCDF file is located;
- **hmin:** Value of *h* parameter;
- **theta\_b:** Control parameter for coordinates close to the ocean floor;
- **theta\_s:** Control parameter for coordinates close to the surface;
- **Tcline:** Critical depth, in meters, controlling the elongation of the coordinates following the terrain. It can be interpreted as the width of the surface where vertical resolution is best;
- **N:** Number of Sigma layers;

- **rmax**: Topography smoothing factor;
- **intrp\_method**: Grid interpolation method: Linear (*linear*) or Near Neighbor (*nn*);
- **grid\_resolution**: Grid resolution;
- **max\_depth**: Maximum grid depth.

**WARNING** In *grid\_resolution*, it is necessary to place the numerator to do the ETOPO1 calculation. For example: as ETOPO1 has  $1/60^\circ$  of spatial resolution, if the chosen spatial resolution is  $1/12^\circ$ , the value of *grid\_resolution* will be 5, because  $5/60^\circ$  is the same as  $1/12^\circ$ .

**WARNING** As we do not use the data from SRTM\_30\_PLUS, the variable *srtm\_dir* does not need to be modified.

```
"""
File name: make_roms_grid.py
Author: Veslei Adriano Sutil
Email: vesleisutil@gmail.com
Created: 17 June 2016
Last modified: 04 July 2019
Version: 6.1

Variable options:
- grd_name      As mentioned in gridid.txt Pyroms file;
- intrp method  'linear' or 'nn' for nearest neighborhood;
- grid_resolution The ETOPO1 spatial resolution is 1/60°. If you desire a 1/12° spatial resolution, enter 5 (5/60 = 1/12°);
The SRTM30+ spatial resolution is 1/120°. If you desire a 1/12° spatial resolution, enter 10 (10/120 = 1/12°).
- max_depth     The grid max depth (m).

Download ETOPO1 here:
https://da.ucar.edu/datasets/ds759.4/index.html#description
Reference:
Amante, C. and B. W. Eakins, 2009: ETOPO1 1 Arc-Minute Global Relief Model: Procedures, Data Sources and Analysis. 24, NOAA Technical Memorandum NESDIS NGDC, 19 pp.

Download SRTM30 plus here:
https://topex.ucsd.edu/WMM/html/srtm30\_plus.html
Reference:
Becker, J. J., D. T. Sandwell, W. H. F. Smith, J. Braud, B. Binder, J. Depner, D. Fabre, J. Factor, S. Ingalls, S-H. Kim, R. Ladner, K. Marks, S. Nelson, A. Pharaoh, R. Trimmer, J. Von Rosenberg, G. Wallace, P. Weatherall., Global Bathymetry

from mpl_toolkits.basemap import Basemap
from bathy_smother import bathy_tools, bathy_smoothing
import mpl_toolkits.basemap as mp
import numpy           as np
import netCDF4
import pyroms
import pyroms_toolbox

# Outputs names.
grd_name      = 'antarctic'
grd_file      = 'grd.nc'
esmpod_dir    = '/home/usuario/model2roms/grid/etopo1.nc'
srtm_dir      = '/home/usuario/model2roms/grid/topo30.nc'

# Grid settings.
hmin          = 25
theta_b       = 2.0
theta_s       = 7.0
tcline         = 200
Tcline         = 40
N              = 40
rmax          = 0.25
intrp_method   = 'linear'
grid_resolution = 5
max_depth     = -5000

# NOTE: Do not need to change above.

# Open bathymetry file.
print("Choose the Digital Elevation Model: (1) ETOPO1 or (2) SRTM30+, then press the [ENTER] button:")
grid_choice = input()
if grid_choice=='1':
    data = netCDF4.Dataset(esmpod_dir, 'r')
    lons = data.variables['x'][:]
    lats = data.variables['y'][:]
    cota = data.variables['z'][:]
    cota = np.array(cota, dtype='float32')
elif grid_choice=='2':
    data = netCDF4.Dataset(srtm_dir, 'r')
    lons = data.variables['x'][:]
    lats = data.variables['y'][:]
    cota = data.variables['z'][:]
    cota = np.array(cota, dtype='float32')
```

**Figure 5.2.** Screenshot of the *make\_grid.py* file, which is located in the *model2roms/grid* folder.

After making the changes, just run the code. Type it:

```
I ipython make_roms_grid.py --pylab
```

### 5.3.3 Introduction to ROMS forcing files

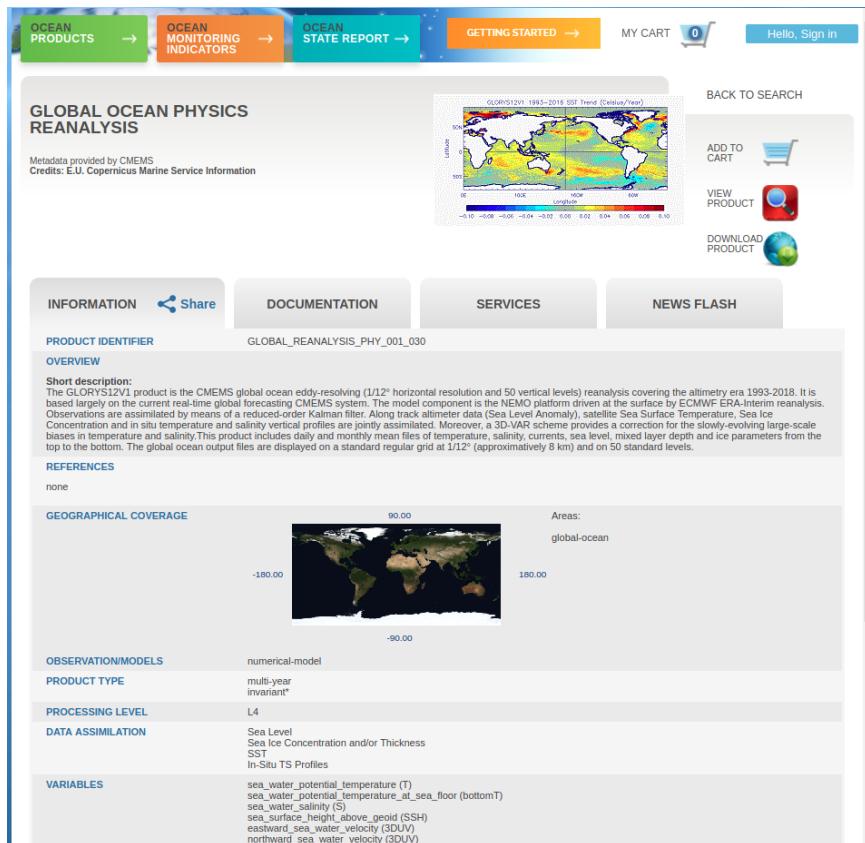
We will use *model2roms* to generate the ROMS conditions. The toolbox supports data from GLORYS12V1 ([Fernandez & Lellouche, 2018](#)) and SODA3 ([Carton et al., 2009](#)), however, we will only use GLORYS12V1 as an example.

### 5.3.4 Global Ocean Physics Reanalysis (GLORYS)

GLORYS is a global reanalysis of the ocean with a spatial resolution of 1/12° with daily or monthly temporal resolution and 50 vertical levels. It is based on the current CMEMS global weather forecasting system and is forced by ECMWF's ERA-Interim from the NEMO model. A Kalman filter with reduced order is used to assimilate the sea altimetry, sea surface temperature, sea ice concentration data obtained from remote sensing and the *in situ* data of vertical temperature and salinity profiles . In addition, a 3D-VAR scheme provides a correction for temperature and salinity deviations.

GLORYS can be downloaded from the website: [http://marine.copernicus.eu/services-portfolio/access-to-products/?option=com\\_csw&view=details&product\\_id=GLOBAL\\_REANALYSIS\\_PHY\\_001\\_030](http://marine.copernicus.eu/services-portfolio/access-to-products/?option=com_csw&view=details&product_id=GLOBAL_REANALYSIS_PHY_001_030).

When you try to download, the site will warn you that you will need to create a Copernicus account. Register and choose the GLORYS daily data set, as shown in Figure 5.3.



**Figure 5.3.** Screenshot from GLORYS webpage.

Click on *Dowload product* and, on the next page (Figure 5.4), select the area according to the latitude and longitude limits of the chosen grid. Also choose the period of data to be downloaded and if you want to use the sea ice model in ROMS, select all variables, with the exception of:

- *bottomT - Sea floor potential temperature;*
- *mlotst - Density ocean mixed layer thickness.*

If you choose to not use the sea ice model, do not select the variables:

- *siconc - Ice concentration;*
- *sithick - Sea ice thickness;*
- *usi - Sea ice eastward velocity;*
- *vsi - Sea ice northward velocity.*

The screenshot shows the 'DATASET FILTERS' section of the GLORYS webpage. At the top, there's a map of the world with green shaded regions indicating 'Intersection between product coverage and area defined by user'. Below the map are two checkboxes: 'Product coverage.' (unchecked) and 'Intersection between product coverage and area defined by user' (checked). A 'Reset geographical selection' button is also present.

**TIME RANGE**  
(Default = Last date available)

START DATE: 2018-12-25 12:00:00  
END DATE: 2018-12-25 12:00:00

**DEPTH**  
(Default = Surface depth)

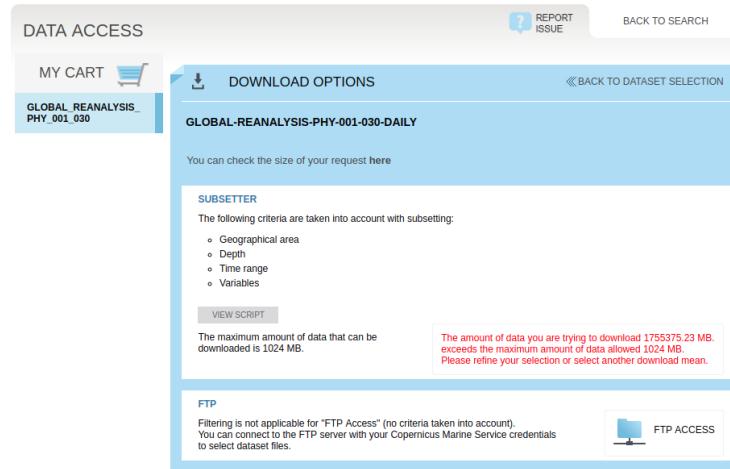
START DEPTH: 0.493  
END DEPTH: 0.4942

**VARIABLES**  
(Default = All variables)

DOWNLOAD	NAME	DESCRIPTION	STANDARD NAME	UNITS
<input checked="" type="checkbox"/>	thetao	Temperature	sea_water_potential_temperature	degrees_C
<input checked="" type="checkbox"/>	bottomT	Sea floor potential temperature	sea_water_potential_temperature_at_sea_floor	degrees_C
<input checked="" type="checkbox"/>	so	Salinity	sea_water_salinity	1e-3
<input checked="" type="checkbox"/>	zos	Sea surface height	sea_surface_height_above_geoid	m
<input checked="" type="checkbox"/>	uo	Eastward velocity	eastward_sea_water_velocity	m s-1
<input checked="" type="checkbox"/>	vo	Northward velocity	northward_sea_water_velocity	m s-1
<input checked="" type="checkbox"/>	mlotst	Density ocean mixed layer thickness	ocean_mixed_layer_thickness_define_d_by_sigma_theta	m
<input checked="" type="checkbox"/>	siconc	Ice concentration	sea_ice_area_fraction	1
<input checked="" type="checkbox"/>	sithick	Sea ice thickness	sea_ice_thickness	m
<input checked="" type="checkbox"/>	usi	Sea ice eastward velocity	eastward_sea_ice_velocity	m s-1

**Figure 5.4.** Another screenshot from the GLORYS webpage.

The Copernicus server allow files for download with a maximum of 1024 mb. If the selected period has a larger file size, as for example in Figure 5.5, it will be necessary to return to the previous step and partition the files into smaller pieces or download the complete file by clicking on *FTP ACCESS* (Figure 5.5).



**Figure 5.5.** Error message when trying to download a text larger than 1024 mb.

If you choose to partition the files or download the files by *FTP*, you will need to create a new file with all the concatenated chosen dates. In this case, it is necessary to use *Climate Data Operators* (CDO).

There are two ways to install CDO: by *Conda* or by *apt-get*.

If you choose to use *Conda*, type:

```
1 conda install -c conda-forge cdo
```

If you choose to use *apt-get*, type:

```
1 sudo apt-get install cdo
```

After installation, enter the directory where all GLORYS files are located and type:

```
1 cdo cat glorys* glorys.nc
```

Where:

- *cdo* means the program command;
- *cat* means the concatenation command of all files;
- *glorys\** means that all files called *glorys* will be concatenated within the folder;

- *glorys.nc* means the final file name created by concatenating with *CDO*.

**WARNING**

Para facilitar os próximos passos, coloque o arquivo do GLORYS dentro do diretório  
\$HOME/model2roms/input

### 5.3.5 Creating ROMS forcing files

When opening the *model2roms* folder, it is possible to observe several script files. We will start with the file *compile.py*, which will compile several files in Fortran90.

Compile the files with the command:

```
1 ipython compile.py --pylab
```

If your *GLORYS* file name is different from *glorys.nc*, open the code *forcingFilenames.py* (Figure 5.6) and in row 15, change '*glorys.nc*' to the name of the created file.

```
1 cd $HOME/model2roms
2 gedit forcingFilenames.py
```

```
import os

# Main function called from model2roms
def getFilename(confM2R, year, month, defaultvar):
    if confM2R.indatatype == 'SODA3':
        if defaultvar is None: defaultvar='salinity'
        filenamein = getSODA3filename(confM2R, year, month, defaultvar)
    if confM2R.indatatype == 'GLORYS':
        if defaultvar is None: defaultvar="temperature"
        filenamein = getGLORYSfilename(confM2R, year, month, defaultvar)
    return filenamein

# private functions called from within module
def getGLORYSfilename(confM2R, year, month, myvar):
    filename = confM2R.modelpath + 'glorys.nc'

    return filename

def getSODA3filename(confM2R, year, month, myvar):
    if (myvar in ['cn', 'hi', 'hs']):
        return confM2R.modelpath + "soda3.3.1_mn_ice_reg_" + str(year) + ".nc"
    else:
        return confM2R.modelpath + "soda3.3.1_mn_ocean_reg_" + str(year) + ".nc"
```

**Figure 5.6.** Screenshot of the *forcingFilename.py* file.

Open the file *configM2R.py* and, in row 16 (Figure 5.7), change the abbreviation *SuaAbreviaçãoAqui* of the *defineabbreviation* to the name of your choice.

```
# Define abbreviation for the run: used to name output files etc.
def defineabbreviation(self):
    return {"SuaAbreviaçãoAqui": "SuaAbreviaçãoAqui"}[self.outgrid]
```

**Figure 5.7.** Screenshot of the *defineabbreviation* in the code *configM2R.py*.

In row 63 (Figure 5.8), change the grid directory *Minha\_Grade.nc* to your grid directory and the abbreviation *SuaAbreviaçãoAqui* by the name chosen in the previous step.

```
def defineromsgridpath(self):
    return {'SuaAbreviaçãoAqui': '/home/usuario/model2roms/grid/Minha_Grade.nc'}[self.outgrid]
```

**Figure 5.8.** Screenshot of the *configM2R.py* code.

On row 67 (Figure 5.9), change the *GLORYS* file directory to your chosen path.

```
def defineforcingdatapath(self):
    return {'SODA3': "/home/usuario/Documentos/model2roms/input/",
           'GLORYS': "/home/usuario/Documentos/model2roms/input/"}[self.indatatype]
```

**Figure 5.9.** Screenshot of the *configM2R.py* code.

From row 76 (Figure 5.10), modify according to your project:

- **self.compileall**: Set as *True* if you want the Fortran files to be recompiled each time *model2roms* is executed;
- **self.createoceanforcing**: Set as *True* to create the hydrodynamic variables;
- **self.createatmosforcing**: Set as *True* to create atmospheric forces. **Currently this function is in the testing phase and is not entirely available for use**;
- **self.writeice**: Set as *True* to create the sea ice variables;
- **self.set2DvarsToZero**: Creates the ice and sea level files with zero values. Since GLORYS has these values, it is recommended to leave it as *False*;
- **self.useesmf**: Set as *True* to use *ESMF* to interpolate the GLORYS data to the ROMS grid;
- **self.usefilter**: Set as *True* to apply a filter to smooth out 2D fields.
- **self.myformat**: The extension for writing ROMS input files. It is, by default, *NETCDF*;
- **self.timefrequencyofinputdata**: The time frequency of the input data. If using GLORYS, write '*day*';
- **self.indatatype**: The name of the data used to generate the forcing files. '*GLORYS* or *SODA3*';
- **self.authorname**: The name of the user using *model2roms*;
- **self.authoremail**: The email of the user using *model2roms*;
- **self.ingridtype**: It will interpolate the GLORYS grid, which is in '*ZLEVEL*' coordinate;
- **self.grdtype**: The GLORYS grid type, which is '*regular*';
- **self.lonname**: The name of the GLORYS longitude variable, which is '*longitude*';
- **self.latname**: The name of the GLORYS latitude variable, which is '*latitude*';
- **self.depthname**: The name of the GLORYS depth variable, which is '*depth*';
- **self.lonname\_u**: The name of the GLORYS U-longitude variable, which is '*longitude*';

- **self.latname\_u**: The name of the GLORYS U-latitude variable, which is '*latitude*';
- **self.lonname\_v**: The name of the GLORYS V-longitude variable, which is '*longitude*';
- **self.latname\_v**: The name of the GLORYS U-latitude variable, which is '*latitude*';
- **self.timename**: The name of the GLORYS time variable, which is '*time*';
- **self.realm**: The realm in which *model2roms* is being executed, in this case, '*ocean*';
- **self.fillvaluein**: The *fillvalue* value of the NetCDF file. By convention, *-1.e20*;
- **self.outgrid**: The name of the abbreviation used in the definition *defineabbreviation*;
- **self.outgridtype**: If "*ROMS*", it will create the output in the standard ROMS format;
- **self.nlevels**: The number of vertical levels in the grid. It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.vstretching**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.vtransform**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.theta\_s**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.theta\_b**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.tcline**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.hc**: It must be the same as the one chosen in the code *make\_roms\_grid.py*;
- **self.start\_year**: The initial year of GLORYS data;
- **self.end\_year**: The final year of GLORYS data;
- **self.start\_month**: The initial month of GLORYS data;
- **self.end\_month**: The final month of GLORYS data;
- **self.start\_day**: The initial day of GLORYS data;
- **self.end\_day**: The final day of GLORYS data;

```

# Set compileAll to True if you want automatic re-compilation of all the
# fortran files necessary to run model2roms. Options are "gfortran" or "ifort". Edit
# compile.py to add other Fortran compilers.
self.compileAll = False
# Create the bry, init, and clim files for a given grid and input data
self.createoceanforcing = True
# Create atmospheric forcing for the given grid
self.createatmosforcing = False # currently in beta stages and unavailable
# Write ice values to file (for Arctic regions)
self.writeice = True
# If you want to require input of lte and ssh, but if you dont have these write zero files to file
self.set2dvarsToZero = False
# Use ESMF for the interpolation. This requires that you have ESMF and ESMPy installed (import ESMF)
self.useesmf = True
# Apply filter to smooth the 2D fields after interpolation (time consuming but enhances results)
self.usefilter = True
# Format to write the output to: 'NETCDF4', 'NETCDF4_CLASSIC', 'NETCDF3_64BIT', or 'NETCDF3_CLASSIC'
# Using NETCDF4 automatically turns on compression of files (ZLIB)
self.myformat = 'NETCDF4'
self.mylib = True
# Frequency of the input data: usually monthly
self.timefrequencyofinputdata = "day" # , "month", "hour"
# Define what grid type you want to interpolate from (input MODEL data)
# Options:
# 1. SIGMA (not properly implemented yet), 2. ZLEVEL
self.ingridtype = "ZLEVEL"
# Define the names of the geographical variables in the input files
self.grdtype = 'regular'
self.lonname = 'longitude'
self.latname = 'latitude'
self.depthname = 'depth'
self.lonname_u = 'longitude'
self.latname_u = 'latitude'
self.lonname_v = 'longitude'
self.latname_v = 'latitude'
self.timename = 'time'
self.realm = 'ocean'
self.levelvaluein = 'e29'
# Define what grid type you want to interpolate to
# Options: This is just the name of your grid used to identify your selection later
self.outgrid = 'antarctic'
self.outgridtype = 'ROMS'
# Define number of output depth levels
self.levels = 40
# Define the grid stretching properties (leave default if uncertain what to pick)
self.vstreching = 4
self.vtransform = 2
self.theta_s = 7.0
self.theta_b = 2.0
self.tcline = 200.0
self.hc = 200
# Define the period to create forcing for
self.start_year = 2015
self.end_year = 2015
self.start_month = 10
self.end_month = 11
self.start_day = 1
self.end_day = 31

```

**Figure 5.10.** Screenshot of the *configM2R.py* code.

Run *model2roms* with the command:

```
! ipython runM2R.py
```

At the end, the initial, boundary and climatological files will be created, which should be added to your project folder, in the Kerana cluster.



## 6. Building the SWAN

SWAN uses the *SWAN.EDT* file to read the grid (*swan\_coord.grd*) and bathymetry (*swan\_bathy.bot*) files. It is also possible to define a numerical grid within the *SWAN.EDT* and indicate the bathymetry file that will be read and associated with the defined grid.

As an example, SWAN files will be created from the ROMS grid, but without the introduction of initial data, as the implementation of a grid inside *SWAN.EDT* is not trivial.

For SWAN boundary conditions, the wind fields will come from the WRF, as soon as the waves generated are associated with the atmospheric systems represented in these simulations. Without the information the boundary files information, the simulated waves will be generated only within the boundary of the domain, disregarding the energy that is leaving or entering the grid.

We will use the MATLAB script, *make\_swan.m*, to generate the files. The script is located at:

```
1 /home/name.surname/repositorio/SWAN_scripts
```

As shown in Figure 6.1, the script has the following construction:

```

clear all
ncfile = '../roms_grid.nc'
x_rho = ncread(ncfile,'lon_rho');
y_rho = ncread(ncfile,'lat_rho');
h = ncread(ncfile,'h');
mask_rho = ncread(ncfile,'mask_rho');

%Replace the land positions with the flag for land (defined in the SWAN
%input file)
land_values = find(mask_rho == 0);
h(land_values) = 9999;

%Print the depths to the bathy file
[n,m] = size(h);

fid = fopen('swan_bathy.bot','w');
for index1 = 1:n;
    for index2 = 1:m;
        fprintf(fid,'   ');
        fprintf(fid,'%12.8f',h(index2,index1));
    end
    fprintf(fid,'\n');
end

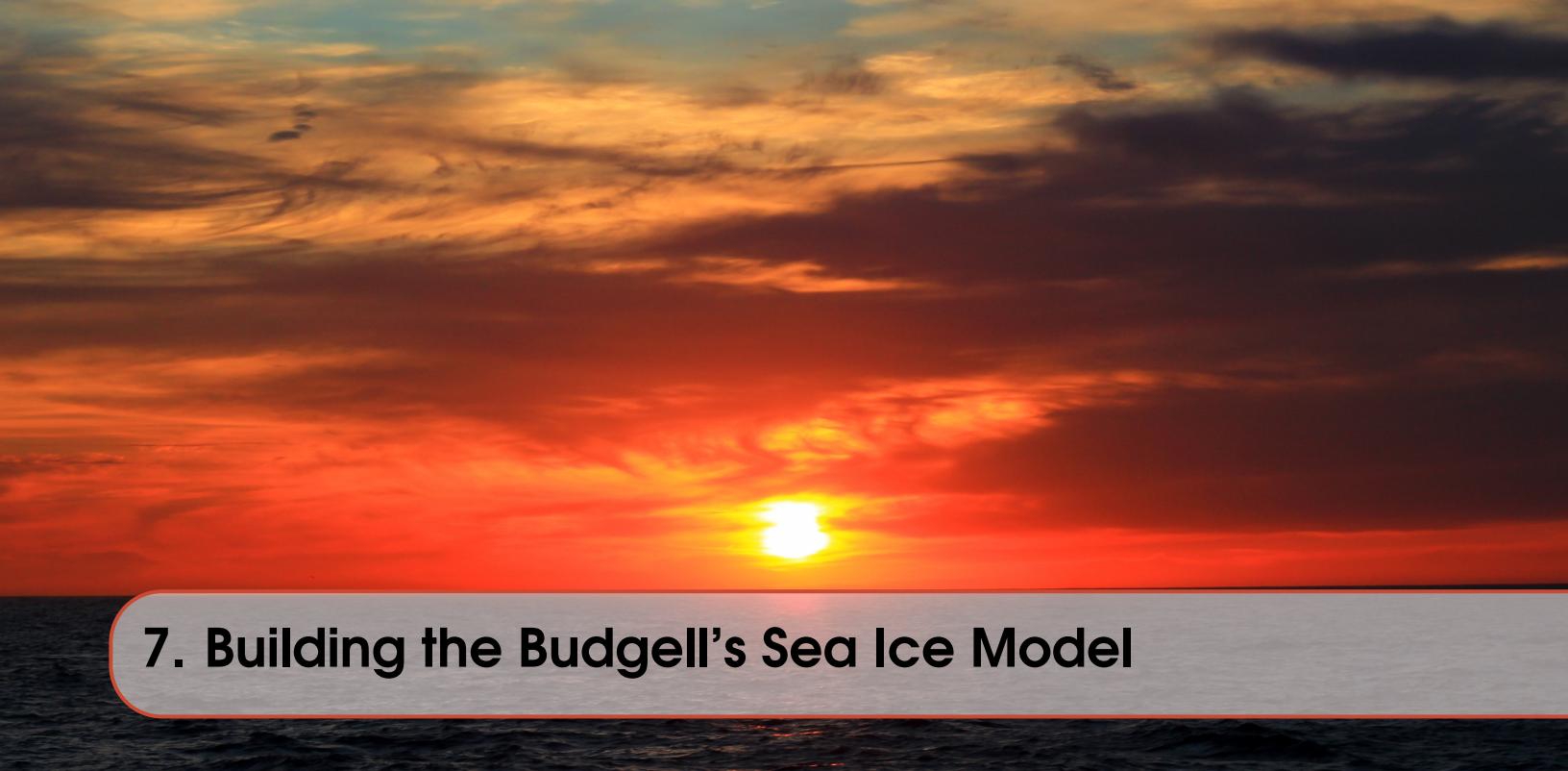
%Print the grid coordinates to the grid file
fid = fopen('swan_coord.grd','w');
fprintf(fid,'%12.6f\n',x_rho);
fprintf(fid,'%12.6f\n',y_rho);

```

**Figure 6.1.** *make\_swam.m* script.

To generate the two SWAN files, look in the script for the variable *ncfile* and change the directory to the path where your ROMS grid is.

Run the script and the two files will be created: *swan\_coord.grd* and *swan\_bathy.bot*. The two files must be placed inside your project folder.



## 7. Building the Budgell's Sea Ice Model

The sea ice model is fully coupled to ROMS, so that, after generating the conditions of the ROMS with ice data (see Section 5.3), to activate the sea ice model just modify the ROMS .h file in your project. According to [Hedström \(2018\)](#), you can add the following options into the .h file, as shown in Figure 7.1:

- **ICE\_MODEL**: defines the sea ice model;
- **ANA\_ICE**: defines initial analytical conditions for sea ice;
- **ICE\_THERMO**: defines the ice thermodynamics;
- **ICE\_MK**: defines the [Mellor & Kanta \(1989\)](#) ice thermodynamics. Currently, is the only choice;
- **ICE\_MOMENTUM**: defines the momentum component of the ice;
- **ICE\_MOM\_BULK**: defines the alternate ice-water stress computation;
- **ICE\_EVP**: defines the elastic-viscous-plastic rheology from [Hunke & Dukowicz \(1997\)](#) and [Hunke \(2001\)](#);
- **ICE\_QUAD\_STRENGTH**: defines the quadratic ice strength from [Overland & Pease \(1988\)](#);
- **ICE\_ADVECT**: defines the advection of ice tracers;
- **ICE\_SMOLAR**: defines the MPDATA use for ice tracers. Currently, it is the only option;
- **ICE\_UPWIND**: defines the upwind advection;
- **ICE\_BULK\_FLUXES**: define the ice part of bulk flux computation;
- **ICE\_DIAGS**: defines the diagnosis of sea ice;
- **ICE\_SHOREFAST** : defines the simple shorefast-ice algorithm from [Budgell \(2005\)](#);
- **ICE\_I\_O**: defines to allow light into the ice as heat;
- **ICE\_CONVSNOW**: defines the conversion of flooded snow to ice.

```
#ifdef SOLVE3D
#define ICE_MODEL
#ifndef ICE_MODEL
#define ANA_ICE
#define ICE_THERMO
#define ICE_MK
#define ICE_MOMENTUM
#define ICE_MOM_BULK
#define ICE_EVP
#define ICE_STRENGTH_QUAD
#define ICE_ADVECT
#define ICE_SMOLAR
#define ICE_UPWIND
#define ICE_BULK_FLUXES
#define ICE_I_O
#define ICE_DIAGS
#endif
#endif
```

**Figure 7.1.** Screenshot of ROMS .h file

For more information about the model, it is recommended to read [Hedström \(2018\)](#).

Get the .in file for the sea ice model. Copy the file *ice.in* into the Kerana repository and add it to your project.

```
1 /home/name.surname/repositorio/ICE_scripts
```

Modify the ROMS .in file, pointing to the *ice.in* into the variable *IPARNAM*:

```
1 nedit ocean.in
2 IPARNAM =  ice.in
```



## 8. Building the weights between grids with SCRIP

### 8.1 Building the weights

As seen in the section 1.6, SCRIP is used to interpolate the weights between two or more grids of different models. In COAWST, the package was modified to generate only one NetCDF file that will be used during integrations.

The SCRIP directory is located at:

```
1 /home/name.surname/COAWST/Lib/SCRIP
```

Inside the folder, look for the file with the extension *.in*. As in the example in Figure 8.1:

```

$INPUTS
| Input file for scrip_coawst
| The $INPUTS line is required at the top of this file.
| Edit this file to enter the correct information below.
| Then run this program as "scrip_coawst scrip_coawst_sandy.in"
| 1) Enter name of output netcdf4 file
OUTPUT_NCFILE='scrip_sandy_moving.nc'
OUTPUT_NCPFILE='scrip_sandy_static.nc'

| 2) Enter total number of ROMS, SWAN, and WRF (max_dom) grids:
NGRIDS_ROMS=2,
NGRIDS_SWAN=2,
NGRIDS_WRF=2,

| 3) Enter name of the ROMS grid file(s):
ROMS_GRIDS(1)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_roms_grid.nc',
ROMS_GRIDS(2)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_roms_grid_ref3.nc',

| 4) Enter SWAN information:
-the name(s) of the SWAN grid file(s) for coords and bathy,
-the size of the SWAN grids, and
-if the swan grids are Spherical(set cartesian=0) or
  Cartesian(set cartesian=1)
SWAN_COORD(1)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_swan_coord.grd',
SWAN_COORD(2)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_swan_coord_ref3.grd',
SWAN_BATH(1)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_swan_bathy.bot',
SWAN_BATH(2)='/home/name.sobrenome/COAWST/Projects/Sandy/Sandy_swan_bathy_ref3.bot',
SWAN_NUMX(1)=84,
SWAN_NUMX(2)=116,
SWAN_NUMY(1)=64,
SWAN_NUMY(2)=86,
CARTESIAN(1)=0,
CARTESIAN(2)=0

| 5) Enter the name of the WRF input grid(s). If the grid is a
moving child nest then enter that grid name as 'moving'.
Also provide the grid ratio, this is used for a moving nest.
WRF_GRIDS(1)='/home/name.sobrenome/COAWST/Projects/Sandy/wrfinput_d01',
WRF_GRIDS(2)='/home/name.sobrenome/COAWST/Projects/Sandy/wrfinput_d02',
WRF_GRIDS(2)='moving',
PARENT_GRID_RATIO(1)=1,
PARENT_GRID_RATIO(2)=3,
PARENT_ID(1)=0,
PARENT_ID(2)=1
PARENT_ID(2)=1

| The $END statement below is required

```

**Figure 8.1.** SCRIP .in file for the Sandy project.

In *OUTPUT\_NCFILE*, change the name of the NetCDF file that will be generated if necessary.

In section 2 of the file, change the variables *NGRIDS\_ROMS*, *NGRIDS\_SWAN* and *NGRIDS\_WRF* according to the number of grids, existing in your project, in ROMS, SWAN and WRF, respectively.

In the third section of the file, renew the ROMS grid directories according to the names in your project.

For SWAN, in the fourth section, in addition to changing the directories of the SWAN grids (*SWAN\_COORD* and *SWAN\_BATH*), change the number of existing grid points, according to your project , in the variables *SWAN\_NUMX* and *SWAN\_NUMY*.

Finally, in the fifth section, change the WRF grid directories (*WRF\_GRIDS*). In *PARENT\_GRID\_RATIO*, if your project includes nesting between the WRF grids, change to the relationship used between the grids used in your project. In *PARENT\_ID*, add the grid ID.

Then, save the changes.

To run SCRIP, search the repository for the file *qsub\_scrip.sh*:

```
| /home/name.surname/repositorio/qsub_scrip.sh
```

Move the file to the SCRIP directory:

```
| mv /home/name.surname/repositorio/qsub_scrip.sh /home/name.surname/COAWST/Lib/SCRIP
```

Open the file *qsub\_scrip.sh*:

```
1 nedit qsub_scrip.sh
```

Change and save the file *.sh*, as in the example in Figure 8.2:

```
#!/bin/sh
#PBS -l mppwidth=1
#PBS -N SCRIP
#PBS -j oe
#PBS -o test.out
#PBS -t 00:00:00-00:00:00
#PBS -q workq
#echo "Running GEOPGRID on KERANA"
#
# export MPICH_ENV_DISPLAY=1
# export MPICH_ABORT_ON_ERROR=1
# export MPICH_RANK_REORDER_DISPLAY=1
# export MPICH_RANK_REORDER_METHOD=1
# export MALLOC_MMAP_MAX_=0
# export OMP_NUM_THREADS=1
# export OMP_NUM_THREADS=1
#
# EXECDIR=/scratch/nome.sobrenome/COAWST/Lib/SCRIP
chmrt
# export ATP_ENABLED=1
# ulimit -a
# ulimit -c unlimited
# ulimit -s unlimited
# ulimit -m unlimited
# ulimit -s

cd $EXECDIR
aprun -n 1 scrip_coawst /scratch/nome.sobrenome/COAWST/Lib/SCRIP/scrip_coawst_sandy.in 1> log.out 2> log.err
```

**Figure 8.2.** *qsub\_scrip.sh* file used to run SCRIP.

To start SCRIP, type:

```
1 qsub qsub_scrip.sh
```

At the end, the file *scrip\_static.nc* will be created. Now put them in your project folder and you're done! COAWST is ready to run.

## 8.2 Running your simulation

Now, with everything ready, your project is ready to be started. Visit the section 2.10 to remember how to execute the project.





## 9. Papers of the Ocean and Atmospheric Studies Laboratory

### Ocean-Atmosphere Interactions in an Extratropical Cyclone in the Southwest Atlantic

U. A. Sutil, L. P. Pezzi, R. C. M. Alves and A. B. Nunes

#### Abstract

This work shows an investigation of the behavior of heat fluxes in the processes of ocean-atmosphere interaction during the passage of an Extra-tropical Cyclone (EC) in the Southwest Atlantic in September 2006 using a coupled regional model's system. A brief evaluation of the simulated data is done by comparison with air and sea surface temperature (SST) data, wind speed, sea level pressure. This comparison showed that both model simulations present some differences (mainly, the wind), nevertheless the simulated variables show quite satisfactory results, therefore allowing a good analysis of the ocean-atmosphere interaction processes. The simulated thermal gradient increases the ocean's heat fluxes into the atmosphere in the cold sector of the cyclone and through the convergence of low level winds the humidity is transported to higher levels producing precipitation. The coupled system showed a greater ability to simulate the intensity and trajectory of the cyclone, compared to the simulation of the atmospheric model.

U. A. Sutil et al. (2018). “Ocean-Atmosphere Interactions in an Extratropical Cyclone in the Southwest Atlantic”. In: *Anuário do Instituto de Geociências - UFRJ*, pp. 525–535. DOI: 10.11137/2019\_1\_525\_535

Available at: [http://www.anuario.igeo.ufrj.br/2019\\_01/2019\\_1\\_525\\_535.pdf](http://www.anuario.igeo.ufrj.br/2019_01/2019_1_525_535.pdf)

**Low connectivity compromises the conservation of reef fishes by marine protected areas in the tropical South Atlantic**

C. A. K. Endo, D. F. M. Gherardi, L. P. Pezzi and L. N. Lima

**Abstract**

The total spatial coverage of Marine Protected Areas (MPAs) within the Brazilian Economic Exclusive Zone (EEZ) has recently achieved the quantitative requirement of the Aichi Biodiversity Target 11. However, the distribution of MPAs in the Brazilian EEZ is still unbalanced regarding the proportion of protected ecosystems, protection goals and management types. Moreover, the demographic connectivity between these MPAs and their effectiveness regarding the maintenance of biodiversity are still not comprehensively understood. An individual-based modeling scheme coupled with a regional hydrodynamic model of the ocean is used to determine the demographic connectivity of reef fishes based on the widespread genus *Sparisoma* found in the oceanic islands and on the Brazilian continental shelf between 10° N and 23° S. Model results indicate that MPAs are highly isolated due to extremely low demographic connectivity. Consequently, low connectivity and the long distances separating MPAs contribute to their isolation. Therefore, the current MPA design falls short of its goal of maintaining the demographic connectivity of *Sparisoma* populations living within these areas. In an extreme scenario in which the MPAs rely solely on protected populations for recruits, it is unlikely that they will be able to effectively contribute to the resilience of these populations or other reef fish species sharing the same dispersal abilities. Results also show that recruitment occurs elsewhere along the continental shelf indicating that the protection of areas larger than the current MPAs would enhance the network, maintain connectivity and contribute to the conservation of reef fishes.

C. A. K. Endo et al. (2019). “Low connectivity compromises the conservation of reef fishes by marine protected areas in the tropical South Atlantic”. In: *Nature Scientific Reports*, pp. 01–11. DOI: 10.1038/s41598-019-45042-0

Available at: <https://www.nature.com/articles/s41598-019-45042-0>

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## An Investigation of Ocean Model Uncertainties Through Ensemble Forecast Experiments in the Southwest Atlantic Ocean

L. N. Lima, L. P. Pezzi, S. G. Penny and C. A. S. Tanajura

### Abstract

Ocean general circulation models even with realistic behavior still incorporate large uncertainties from external forcing. This study involves the realization of ensemble experiments using a regional model configured for the Southwest Atlantic Ocean to investigate uncertainties derived from the external forcing such as the atmosphere and bathymetry. The investigation is based on perturbing atmospheric surface fluxes and bathymetry through a series of ensemble experiments. The results showed a strong influence of the South Atlantic Convergence Zone on the underlying ocean, 7 days after initialization. In this ocean region, precipitation and radiation flux perturbations notably impacted the sea surface salinity and sea surface temperature, by producing values of ensemble spread that exceeded 0.08 and 0.2 °C, respectively. Wind perturbations extended the impact on currents at surface, with the spread exceeding 0.1 m/s. The ocean responded faster to the bathymetric perturbations especially in shallow waters, where the dynamics are largely dominated by barotropic processes. Ensemble spread was the largest within the thermocline layer and in ocean frontal regions after a few months, but by this time, the impact on the modeled ocean obtained from either atmospheric or bathymetric perturbations was quite similar, with the internal dynamics dominating over time. In the vertical, the sea surface temperature exhibited high correlation with the subsurface temperature of the shallowest model levels within the mixed layer. Horizontal error correlations exhibited strong flow dependence at specific points on the Brazil and Malvinas Currents. This analysis will be the basis for future experiments using ensemble-based data assimilation in the Southwest Atlantic Ocean.

L. N. Lima et al. (2019). “An Investigation of Ocean Model Uncertainties Through Ensemble Forecast Experiments in the Southwest Atlantic Ocean”. In: *Journal of Geophysical Research: Oceans* 120, pp. 432–452

Available at: <https://agupubs.onlinelibrary.wiley.com/doi/epdf/10.1029/2018JC013919>

**Coupled ocean-atmosphere forecasting at short and medium time scales**

J. Pullen, R. Allard, H. Seo, A. J. Miller, S. Chen, L. P. Pezzi, T. Smith, P. Chu, J. Alves and R. Caldeira

**Abstract**

Recent technological advances over the past few decades have enabled the development of fully coupled atmosphere-ocean modeling prediction systems which are used today to support short-term (days to weeks) and medium-term (10-21 days) needs for both the operational and research communities. Utilizing several coupled modeling systems we overview the coupling framework, including model components and grid resolution considerations, as well as the coupling physics by examining heat fluxes between atmosphere and ocean, momentum transfer, and freshwater fluxes. These modeling systems can be run as fully coupled atmosphere-ocean and atmosphere-ocean-wave configurations. Examples of several modeling systems applied to complex coastal regions including Madeira Island, Adriatic Sea, Coastal California, Gulf of Mexico, Brazil, and the Maritime Continent are presented. In many of these studies, a variety of field campaigns have contributed to a better understanding of the underlying physics associated with the atmosphere-ocean feedbacks. Examples of improvements in predictive skill when run in coupled mode versus standalone are shown. Coupled model challenges such as model initialization, data assimilation, and earth system prediction are discussed.

J. Pullen et al. (2017). *The Science of Ocean Prediction, The Sea*. P. Lermusiaux and K. Brink. Chap. Coupled ocean-atmosphere modeling and predictions

Available at: [http://meteora.ucsd.edu/~miller/papers/TheSea\\_Chapter23.html](http://meteora.ucsd.edu/~miller/papers/TheSea_Chapter23.html)

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## Regional modeling of the water masses and circulation annual variability at the Southern Brazilian Continental Shelf

L. F. Mendonça, R. B. Souza, C. R. C. Aseff, L. P. Pezzi, O. O. Möller and R. C. M. Alves

### Abstract

The Southern Brazilian Continental Shelf (SBCS) is one of the more productive areas for fisheries in Brazilian waters. The water masses and the dynamical processes of the region present a very seasonal behavior that imprint strong effects in the ecosystem and the weather of the area and its vicinity. This paper makes use of the Regional Ocean Modeling System (ROMS) for studying the water mass distribution and circulation variability in the SBCS during the year of 2012. Model outputs were compared to in situ, historical observations and to satellite data. The model was able to reproduce the main thermohaline characteristics of the waters dominating the SBCS and the adjacent region. The mixing between the Subantarctic Shelf Water and the Subtropical Shelf Water, known as the Subtropical Shelf Front (STSF), presented a clear seasonal change in volume. As a consequence of the mixing and of the seasonal oscillation of the STSF position, the stability of the water column inside the SBCS also changes seasonally. Current velocities and associated transports estimated for the Brazil Current (BC) and for the Brazilian Coastal Current (BCC) agree with previous measurements and estimates, stressing the fact that the opposite flow of the BCC occurring during winter in the study region is about 2 orders of magnitude smaller than that of the BC. Seasonal maps of simulated Mean Kinetic Energy and Eddy Kinetic Energy demonstrate the known behavior of the BC and stressed the importance of the mean coastal flow off Argentina throughout the year.

L. F. Mendonça et al. (2017). “Regional modeling of the water masses and circulation annual variability at the Southern Brazilian Continental Shelf”. In: *Journal of Geophysical Research: Oceans* 122, pp. 1232–1253.  
DOI: 10.1002/2016JC011780

Available at: <https://agupubs.onlinelibrary.wiley.com/doi/abs/10.1002/2016JC011780>

**The Influence of Sea Ice Dynamics on the Climate Sensitivity and Memory to Increased Antarctic Sea Ice**

C. K. Parise, L. P. Pezzi, K. I. Hodges and F. Justino

**Abstract**

The study analyzes the sensitivity and memory of the Southern Hemisphere coupled climate system to increased Antarctic sea ice (ASI), taking into account the persistence of the sea ice maxima in the current climate. The mechanisms involved in restoring the climate balance under two sets of experiments, which differ in regard to their sea ice models, are discussed. The experiments are perturbed with extremes of ASI and integrated for 10 yr in a large 30-member ensemble. The results show that an ASI maximum is able to persist for 4 yr in the current climate, followed by a negative sea ice phase. The sea ice insulating effect during the positive phase reduces heat fluxes south of 60°S, while at the same time these are intensified at the sea ice edge. The increased air stability over the sea ice field strengthens the polar cell while the baroclinicity increases at midlatitudes. The mean sea level pressure is reduced (increased) over high latitudes (midlatitudes), typical of the southern annular mode (SAM) positive phase. The Southern Ocean (SO) becomes colder and fresher as the sea ice melts mainly through sea ice lateral melting, the consequence of which is an increase in the ocean stability by buoyancy and mixing changes. The climate sensitivity is triggered by the sea ice insulating process and the resulting freshwater pulse (fast response), while the climate equilibrium is restored by the heat stored in the SO subsurface layers (long response). It is concluded that the time needed for the ASI anomaly to be dissipated and/or melted is shortened by the sea ice dynamical processes.

C. K. Parise et al. (2014). “The Influence of Sea Ice Dynamics on the Climate Sensitivity and Memory to Increased Antarctic Sea Ice”. In: *Journal of Climate* 28, pp. 9642–9668. DOI: 10.1175/JCLI-D-14-00748.1

Available at: <https://journals.ametsoc.org/doi/10.1175/JCLI-D-14-00748.1>

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**Modeling the spawning strategies and larval survival of the Brazilian sardine (*Sardinella brasiliensis*)**

D. F. Dias, L. P. Pezzi, D. F. M. Gherardi and R. Camargo

**Abstract**

An Individual Based Model (IBM), coupled with a hydrodynamic model (ROMS), was used to investigate the spawning strategies and larval survival of the Brazilian Sardine in the South Brazil Bight (SBB). ROMS solutions were compared with satellite and field data to assess their representation of the physical environment. Two spawning experiments were performed for the summer along six years, coincident with ichthyoplankton survey cruises. In the first one, eggs were released in spawning habitats inferred from a spatial model. The second experiment simulated a random spawning to test the null hypothesis that there are no preferred spawning sites. Releasing eggs in the predefined spawning habitats increases larval survival, suggesting that the central-southern part of the SBB is more suitable for larvae development because of its thermodynamic characteristics. The Brazilian sardine is also capable of exploring suitable areas for spawning, according to the interannual variability of the SBB. The influence of water temperature, the presence of Cape Frio upwelling, and surface circulation on the spawning process was tested. The Cape Frio upwelling plays an important role in the modulation of Brazilian sardine spawning zones over SBB because of its lower than average water temperature. This has a direct influence on larval survival and on the interannual variability of the Brazilian sardine spawning process. The hydrodynamic condition is crucial in determining the central-southern part of SBB as the most suitable place for spawning because it enhances simulated coastal retention of larvae.

D. F. Dias et al. (2014). “Modeling the spawning strategies and larval survival of the Brazilian sardine (*Sardinella brasiliensis*)”. In: *Progress in Oceanography* 123, pp. 38–53. doi: 10.1016/j.pocean.2014.03.009

Available at: <http://www.iag.usp.br/pos/meteorologia/biblio/modeling-spawning-strategies-and-larval-survival-brazilian-sardine-sardinella-br>

**Sea surface temperature anomalies driven by oceanic local forcing in the Brazil-Malvinas Confluence**

I. P. da Silveira and L. P. Pezzi

**Abstract**

Sea surface temperature (SST) anomaly events in the Brazil-Malvinas Confluence (BMC) were investigated through wavelet analysis and numerical modeling. Wavelet analysis was applied to recognize the main spectral signals of SST anomaly events in the BMC and in the Drake Passage as a first attempt to link middle and high latitudes. The numerical modeling approach was used to clarify the local oceanic dynamics that drive these anomalies. Wavelet analysis pointed to the 8–12-year band as the most energetic band representing remote forcing between high to middle latitudes. Other frequencies observed in the BMC wavelet analysis indicate that part of its variability could also be forced by low-latitude events, such as El Niño. Numerical experiments carried out for the years of 1964 and 1992 (cold and warm El Niño-Southern Oscillation (ENSO) phases) revealed two distinct behaviors that produced negative and positive sea surface temperature anomalies on the BMC region. The first behavior is caused by northward cold flow, Río de la Plata runoff, and upwelling processes. The second behavior is driven by a southward excursion of the Brazil Current (BC) front, alterations in Río de la Plata discharge rates, and most likely by air-sea interactions. Both episodes are characterized by uncoupled behavior between the surface and deeper layers.

I. P. Silveira & L. P. Pezzi (2014). “Sea surface temperature anomalies driven by oceanic local forcing in the Brazil-Malvinas Confluence”. In: *Ocean Dynamics* 347.64, pp. 347–360. DOI: 10.1007/s10236-014-0699-4

Available at: <https://link.springer.com/article/10.1007/s10236-014-0699-4>



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