

Activity 4 – Set up and run a realistic configuration

0. If you need to connect to LOPS computers:

ssh -Y username@polaris.univ-brest.fr

ssh -Y libra or *ssh -Y capella* or *ssh -Y apus*

cd /net/krypton/data0/project/vortex/gula/ModNum/username (ask me if it is not created)

1. Download the croco tools and datasets

- copy the CROCO_TOOLS package from:

if you are logged in a LOPS computer:

/net/krypton/data0/project/vortex/gula/ModNum/Data/croco_tools-v1.1.tar.gz

if you are in a IUEM computing rooms:

/forums/public/pub/Data/ croco_tools-v1.1.tar.gz

Otherwise you need to download it from

<https://www.croco-ocean.org/download/croco-project/>

or

http://jgula.fr/ModNum/croco_tools-v1.1.tar.gz

- We are going to use the Matlab version. Untar the folder (*tar xf croco_tools-v1.1.tar.gz*) and copy the files *croco_tools-v1.1/start.m* and *croco_tools-v1.1/crocotools_param.m* in your working folder (e.g. *~/ModNum/myconfig*)

- edit the *start.m* and modify the path to the croco_tools routines if needed (tools_path needs to point to the *croco_tools-v1.1* folder)

- Edit the `crocotools_param.m` to and modify the path to the data files (variable DATADIR). It should point to:

if you are logged in a LOPS computer:

`/net/krypton/data0/project/vortex/gula/ModNum/Data/`

if you are in a IUEM computing rooms:

`/forums/public/pub/Data/`

Otherwise you need to download datasets (COADS + WOA + Topo) from

<https://www.croco-ocean.org/download/datasets/>

2. Create files for the Benguela test-case

See https://croco-ocean.gitlabpages.inria.fr/croco_doc/ for a detailed tutorial

In brief you need to :

- start matlab:

On LOPS computers: ***module load matlab/2014a; matlab -nodesktop***

and run `start.m` to load all path (*just type "start" in the matlab window/terminal*)

- run the following scripts (just type their name in matlab)

- `make_grid` to create your grid file [CROCO_FILES/croco_grd.nc]
- `make_forcing` to generate surface_forcing: wind stress, surface heat flux, surface freshwater flux [CROCO_FILES/croco_frc.nc]
- `make_clim` to generate initial conditions T, S, currents , SSH [CROCO_FILES/croco_ini.nc and CROCO_FILES/croco_clm.nc]
- `make_bry` to generate oceanic boundary conditions : T, S, currents , SSH [CROCO_FILES/croco_bry.nc]

3. Compile and run the BENGUELA_LR case

Redownload and untar the code croco if you don't have it:

```
cd .. (wherever you want to have your croco routines)  
wget https://www.jgula.fr/ModNum/croco.tar.gz  
tar xf croco.tar.gz  
cd regional (the folder containing your latest configuration)
```

- Copy the jobcomp locally: **cp -rf ../croco/OCEAN/jobcomp ./**
- Copy the croco.in locally: **cp -rf ../croco/OCEAN/croco.in ./**
- Copy the param.h locally: **cp -rf ../croco/OCEAN/param.h ./**
- Copy the cppdefs.h locally: **cp -rf ../croco/OCEAN/cppdefs.h ./** or **wget <http://jgula.fr/ModNum/cppdefs.h>** if you have the previous croco version

On LOPS computers:

- Load the fortran/netcdf modules:
module purge
module load intel/12.1 netcdf/c-4.4.1.1-intel12 netcdf/fortran-4.4.4-intel12
- and add the corresponding locations in the jobcomp:
NETCDFINC="-I/opt/linux/netcdf/fortran-4.4.4-intel12/include -I/opt/linux/netcdf/4.4.1.1-intel12/include"
NETCDFLIB="-L/opt/linux/netcdf/fortran-4.4.4-intel12/lib -L/opt/linux/netcdf/4.4.1.1-intel12/lib -lnetcdf -lnetcdff"
- eventually change the SOURCE location in the jobcomp to point to the croco routines:
SOURCE=../croco/OCEAN

Add OPEN_MP parallelization (see Activity1.pdf)

Then Compile and run :

./jobcomp
croco croco.in

4. Create files for your own configuration

- edit the [crocotools_param.m](#) to choose all parameters for your configuration (name, grid location and size, time and duration, path to forcing files, etc.)

5. Compile and run your simulation

- Edit the [param.h](#) to define the size of your grid (and parameters for your parallelization) according to the ones you have chosen in [crocotools_param.m](#)
- Edit the [cppdefs.h](#) to choose your numerical options
- Edit the [croco.in](#) to choose the run time parameters. Choose the parameters to have at least 2 years of simulation with monthly averages (every 30 days).

6. Check your simulation

- Check the circulation:

- you can use a [python gui](#) in python, see https://croco-ocean.gitlabpages.inria.fr/croco_doc/tutos/tutos.14.visu.python.html

- you can use the [croco_gui.m](#) in with matlab, see https://croco-ocean.gitlabpages.inria.fr/croco_doc/tutos/tutos.14.visu.matlab.html

- or various examples in python (see [Modules.zip](#), [Python_example1.py](#), [Python_example2.py](#) on <http://jgula.fr/ModNum>)

- Try to find a paper documenting the circulation in this region and check if it is (at least qualitatively) well reproduced in the simulation.

- Average the simulation over the last year only (*you will consider only the last year of your simulation to minimize the effects of the spin-up.*)
- Plot the mean currents (surface and barotropic) and vertical sections of stratification for your simulation and for observations (*you can use WOA2009 data – or directly use the croco_clm.nc file which contains monthly climatology from WOA2009 data interpolated on the model grid*).