Activity 1 - Run an idealized ocean gyre

1. Get the model

- Download from http://jgula.fr/ModNum/croco.tar.gz or from /forums/public/pub/Data/croco.tar.gz
- Put it in a folder of your choice: [for example ~/ModNum/]
- untar it: tar xf croco.tar.gz
- the croco source code is in the folder ~/ModNum/croco/OCEAN

2. Compile the model

- Create a folder where you will run the model [for example ~/ModNum/case1]
- we need to edit the following files: jobcomp, cppdefs.h, param.h, croco.in so copy them into the folder you just created:

```
cp ~/ModNum/croco/OCEAN/jobcomp ~/ModNum/case1/
```

- cp ~/ModNum/croco/OCEAN/cppdefs.h ~/ModNum/case1/
- cp ~/ModNum/croco/OCEAN/param.h ~/ModNum/case1/
- cp ~/ModNum/croco/TEST_CASES/croco.in.Basin ~/ModNum/case1/croco.in
- the model needs a fortran compiler and compatible netcdf libraries:
 - If you are working with Linux on IUEM computers:
 - ★1. unload everything: module purge
 - ★2. then load intel compilers and netcdf library module load intel/12.1 netcdf/4.1.3-intel12
 - If you are working with MacOS:
 - * you need to have gcc and netcdf installed (using homebrew or macports): brew install gcc; brew install netcdf
 - *you may need to edit the jobcomp to specify where to find the netcdf libraries (nf-config doesn't work with homebrew version of netcdf-fortran)

NETCDFLIB="-L/usr/local/lib -lnetcdff" NETCDFINC="-I/usr/local/include" #NETCDFLIB=\$(nf-config --flibs) #NETCDFINC=-I\$(nf-config --includedir) - edit the jobcomp to specify the location of the source code:

SOURCE=../croco/OCEAN

- edit the cppdefs.h and choose the predefined test case **Basin**:

#define BASIN
...
#undef REGIONAL

- Compile the code: ./jobcomp

3. Run the model

days)

- run the model: ./croco croco.in
- look at the output: neview basin_his.ne

4. Modify the namelist (croco.in)

- Modify the namelist (croco.in) to:
 - a. make the model run for 20 years (you can approximate 1 year = 360
 - b. output of history files every 30 days
 - c. outputs of variables averaged over 5 years

- Note that you also need to modify the cppdefs.h to output averaged files:

5. Run the model using openMP

- we need to edit the file cppdefs.h,
- find the part of the file where the BASIN case is defined (#elif defined BASIN) and change # undef OPENMP into # define OPENMP

- edit the param.h file to choose the number of processors (NPP) you want to use:

```
#elif defined OPENMP
    parameter (NPP=8)
```

- reCompile the code: ./jobcomp
- define the number of threads to use for your computer (the number should be the same than NPP defined in the param.h file), type in your terminal:
 - in bash: export OMP_NUM_THREADS=8in csh/tcsh: setenv OMP_NUM_THREADS 8
- run the model: ./croco croco.in
- What can you tell about the configuration? What is the forcing? What is the Coriolis parameter? What is the boundary condition? What is the bottom condition?

- What can you tell about the circulation? Is it similar to Stommel's gyre? (http://jgula.fr/ModNum/Stommel48.pdf)