
CLIMATE MODELLING DIAGNOSTICS

SOFTWARE: **EZCLIM2.0**

USER GUIDE

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IF YOU COME ACROSS ANY PROBLEMS, SEE [SECTION 8](#) FOR POSSIBLE
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Getting started with the Climate Modelling Software

The Climate Modelling (Diagnostics) Software is written using Python 3 and enables users to compute climate model diagnostics. You can run it either on JASMIN or Locally follow these steps.

1.1 Quick setup

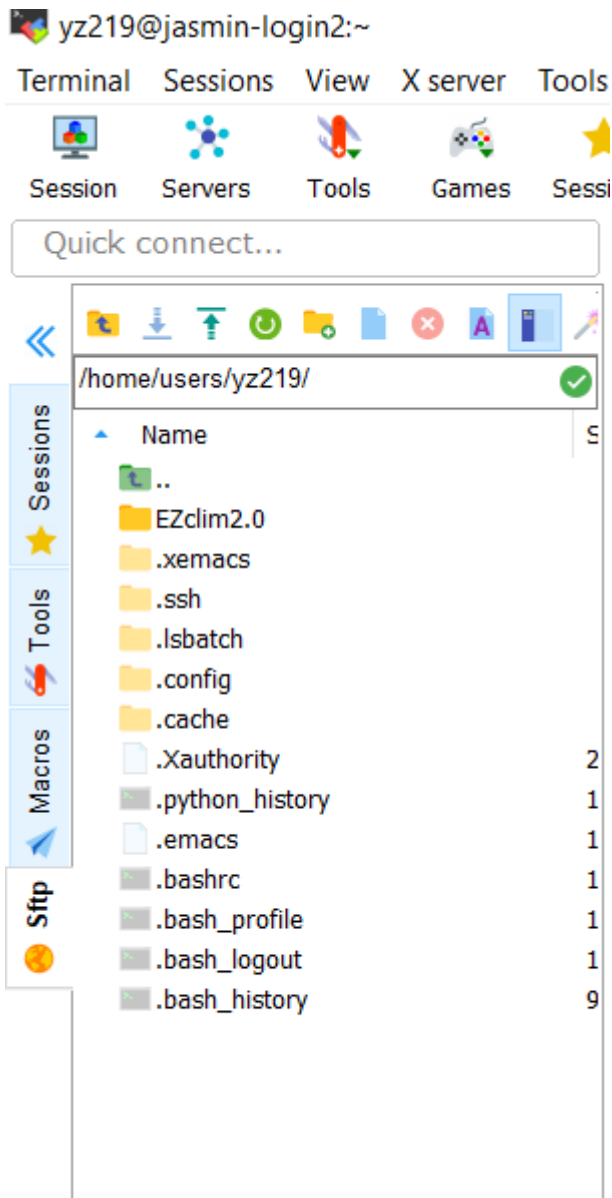
These steps give a quick setup for running on JASMIN. If you get stuck on any of the steps or you want to run it locally, please start from section 1.2.

- Download / clone the repository into your computer from here (<https://github.com/acse-yz219/EZclim2.0>)
- You must first apply for a JASMIN Account (<https://accounts.jasmin.ac.uk>)
- Once your JASMIN Account is successfully created, you need to apply for a JASMIN Service (https://accounts.jasmin.ac.uk/services/my_services/)
- Once you have the permission to access JASMIN Service, download MobaXterm, (<https://mobaxterm.mobatek.net/>)
MobaXterm is the recommended ssh server of JASMIN to connect its service.

- Go into MobaXterm:

Follow the instruction of (<https://help.jasmin.ac.uk/article/187-login>) to login to the JASMIN service

- Once you can successfully login to the JASMIN Service, upload the whole EZclim2.0 to the Service. Just drag the file into the box.



- Run module load jasper to enable python3
- cd EZclim2.0

1.2 Downloading package

The software package is stored in [GitHub](#), a repository hosting service. You can either download the folder **manually** or through the Linux **terminal**.

1.2.1 Manual download

1. Go to the the repository linked [here](https://github.com/acse-yz219/EZclim2.0). (<https://github.com/acse-yz219/EZclim2.0>)
2. Go the upper right hand corner and click on the **Clone or download** button. Then click on **Download Zip**.

3. Extract the folder from the zipped file that has downloaded.
4. Go to a command line terminal and go to the directory where the folder is stored.
Then:

```
$ cd EZclim2.0
```

5. To check the contents in the folder, do:

```
$ ls
```

If you do not see the folder **EZclim2.0** in the list, then you may need to **cd** again.

If you can see the folders **EZclim2.0**, then the package has successfully downloaded.

1.2.2 Terminal download (Suitable for Linux and Mac)

1. Go to the terminal and check that since we are installing through *Git*, we have to ensure that it is installed. Type on the command line:

```
$ git --version
```

If this gives an output like git version 2.17.1 (the numbers can be different), then git is already installed.

If git is not installed, then to install type on the command line:

```
$ sudo apt install git-all
```

Note: If **sudo** does not work, remove it and just put apt install git-all. After it has finished installing, check for the version (as in the previous step) to make sure it has installed properly.

2. To download from git, you have to clone the repository (the link has no spaces):

```
$ git clone https://github.com/acse-yz219/EZclim2.0.git
```

3. Go into the folder to make sure it is cloned properly

```
$ cd EZclim2.0
```

and

```
$ ls
```

If you can see the folders **EZclim2.0**, then the package has successfully downloaded.

1.3 Installing all requirements (do only if you want to run it locally)

The following instructions are suitable for Linux, macOS or Windows (with an Ubuntu subsystem or Anaconda).

1. First ensure that Python is installed on the machine.

```
$ python --version
```

If you can see a version number, then Python is installed.

If Python is not installed, then install it by:

```
$ sudo apt-get update  
$ sudo apt-get install python3.6
```

2. You need to install Miniconda in order to get some of the libraries. (The link has no spaces).

```
$ wget https://repo.continuum.io/miniconda/Miniconda3-latest-  
Linux-x86_64.sh
```

```
$ chmod +x Miniconda3-latest-Linux-x86_64.sh
```

```
$ ./Miniconda3-latest-Linux-x86_64.sh
```

After this you should see a "Welcome to Miniconda", and pressing Enter, you will see a license to accept.

Press enter until you get to the end of the licence and type `yes` to accept.

Press enter to confirm the location. Type `yes` for the installer to initialise Miniconda by running `conda init`.

After packages are installed, you should see a message saying "Thank you for installing Miniconda3".

Restart the terminal so that all the changes can take effect.

After restarting, check conda has installed by checking the version:

```
$ conda --version
```

3. We can now install the rest:

```
$ conda install -c conda-forge iris
```

And for the plotting (used to overlay the continents on the map plots):

```
$ conda install basemap
```

4. To install the rest of the dependencies, make sure you are in the git repo folder `acse-9-independent-research-project-AdannaAkwats` folder and see the **EZ-clim** and **EZclim_Parallel** folders. Then do:

```
$ pip install -r INSTALL.txt
```

At the end, it should have: "Successfully installed Cython-0.29.13 Shapely-1.6.4.post2 attrs-19.1.0 cdo-1.5.3 more-itertools-7.2.0 netCDF4-1.4.2 numpy-1.16.3 pluggy-0.6.0 py-1.8.0 pytest-3.5.1 python-dateutil-2.7.3 scipy-1.2.1" seaborn-0.9.0 xarray-0.12.3" or something similar.

5. If pip is not installed, then install pip with:

```
$ sudo apt-get python -pip
```

and then go back to step 4. As before, if using `sudo` does not work, just use `apt-get python-pip` instead.

We now have all packages installed!

Some example code is available in [section 3](#) to run.

1.4 Installing all requirements (for Windows without Anaconda or Ubuntu subsystem)(do only if you want to run it locally)

In Windows, there are 2 ways to run Python programs - either by downloading **Anaconda** (a Python Data Science Platform) or setting up and downloading the **Windows Subsystem for Linux**.

1.4.1 Using Anaconda

Follow these steps on the official Anaconda site to download Anaconda for Windows:

<https://docs.anaconda.com/anaconda/install/windows/>. Download the **Python 3.7 version**.

After set up is complete, open the Anaconda command prompt, go back to step 2 in [section 1.3](#).

1.4.2 Using the Windows Ubuntu subsystem

Follow the steps given in the official Microsoft web page <https://docs.microsoft.com/en-us/windows/wsl/install-win10>. I recommend downloading one of the **Ubuntu** terminals.

After set up is complete, open the Ubuntu terminal, go back to step 1 in [section 1.3](#).

Package structure

The package is structured as follows.

EZclim 2.0

|--

|-- EZclim/

| |-- DATA/

| |-- INPUT/

| |-- user_function/

| |-- input.txt

| |-- input_example.txt

| |-- Preprocessing.txt

|

| |-- RESULTS/

| |-- ensemble_averages/

|

| |-- directories.py

| |-- output.log

| |-- main.py

|

|-- INSTALL.txt |-

- README

|-- UserGuide.pdf

- DATA/ : The default folder where NetCDF model output files are stored. You can redirect this directory in `directories.py` if the data is stored somewhere else.
- INPUT/ : Folder where all input files (e.g `input.text`, mask files etc) are stored.
- RESULTS/ : Folder where all results and files from software run is stored. In `ensemble_averages`, the NetCDF output files from program are stored.
- `output.log` : File where all progress and error messages are stored.



Running the software

3.1 Preprocessing

Enter the **EZclim2.0** folder on JASMIN:

```
$ cd EZclim2.0
```

Filling the preprocessing.txt and run

```
$ python Preprocessing.py
```

It usually takes a long time, be patient. If there are some warnings, don't worry. This is because in re-grid stage, there are many null values in the original file.

3.2 Diagnostics

Enter the **EZclim** folder on JASMIN:

```
$ cd EZclim2.0
```

There are 2 ways to run the diagnostics, either using an input file or using command line arguments.

1. To run the input_example.txt file which calculates the mean of air temperature in North America and plots the map and histogram:

```
$ python main.py -ex
```

After running, it should show the plots and print a message on the command line when plots are closed.

NOTE: If you run this on a terminal does not allow plotting, an error will be thrown. So you have to set plot to "False" in the input file or go to the [Troubleshooting section](#).

2. After filling the arguments in input.txt, the code can be run with this input by:

```
$ python main.py
```

3. You can also use the command line, e.g. to run with the input_example.txt arguments:

```
$ python main.py 1970 -pf E1_north_america -v air_temperature -e 1  
-a mean -p 1 -o
```

For help on what each argument means:

```
$ python main.py --help
```

4. After each run, you can check the progress or potential error messages by:

```
$ less output.log
```

3.2.1 Debugging

When the program is run successfully, the following message prints out to the terminal:

PROGRAM FINISHED.

>> Progress and potential errors are logged in output.log file.

>> To open output.log, type in cmd line: less output.log

If no message prints out, it is because an error has been seen in the code, this may be caused by invalid inputs being seen. Open the output.log file created and this will give you a summary of what the code has been doing, and any errors that may have been thrown.

4

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Software input

This section details the input required by the package.

4.1 File format

The names of the files in the default DATA folder should be in the general format. The file on JASMIN already meet the requirement, but the output file name should still meet the requirement.

"{START OF FILENAME}_{YEAR}.nc"

where

- {START OF FILENAME} is the prefix of the file
- {YEAR} is the year

Note that the separation underscore ' _ ' before the year does not need to be an underscore. E.g. it can be separated by a dot.

You can also add the *ensemble number* in the file name:

"{START OF FILENAME}_ens{NUM}_{YEAR}.nc"

where

- {START OF FILENAME} is the prefix of the file
- {YEAR} is the year
- {NUM} is the ensemble number

Value of {NUM}: If the number of ensembles are less than 100, then the values of {NUM} should be 101, 102, ..., 199. This represent the 1st, 2nd ...99th ensemble member. The '1' given at the start is given for readability. Similarly if the number of ensembles is less than 1000, then values of {NUM} are 1001, 1002,1100, 1102, ...1999, for the 1st all the way to the 999th member.

If the file name does not have the ens{NUM} then it is assumed that the model output is the first (and only) ensemble member.

For example, if I had 3 ensembles of the model output for air temperature at 1970. I could have the file names:

- air_temp_ens101_1970.nc
- air_temp_ens102_1970.nc
- air_temp_ens103_1970.nc

If you have multiple years stored in one file then the format is:

"{START OF FILENAME}_{YEAR1}_{YEAR 2}.nc"

or

"{START OF FILENAME}_ens{NUM}_{YEAR1}-{YEAR 2}.nc"

where

- {START OF FILENAME} is the prefix of the file
- {YEAR 1} and {YEAR 2} are the start and end year of the data in the file.

Example file names:

Some valid example names are:

- dic_deltap_ens102_1d_1953.nc
- cCwd_ens101_2015_2065.nc
- pcmdi.ipcc4.mri_cgcm2_3_2a.picntrl.run1.annual.thetao_O1.1990s.nc

(Files from CMIP6):

- thetaoga_Omon_EC-Earth3-Veg_historical_r1i1p1f1_gn_185301-185312.nc
- ccb_Amon_CNRM-CM6-1_highresSST-present_r1i1p1f2_gr_200001-200912.nc

Invalid example names:

- dic_deltap.nc

A year **must** be added to the file name.

4.2 File data assumptions

- Multiple files do not have overlapped data. (i.e. 2 files should not have the same data)

- Daily or monthly increments of data, higher resolution of data input e.g. hourly is not supported.
- The time increments of data must be consistent. E.g. if you are passing 2 NetCDF files as input, **both** must have either monthly or daily data.
- Only rectangular grids are supported. If (for example) the model uses a curvilinear grid, this will have to be converted to a rectangular grid before passing into the program.

4.3 Input arguments for Preprocessing

The following are the required arguments needed in the Preprocessing

4.3.1 File Path

File Path: File Path is the file path of the file that you want to process on JASMIN

4.3.2 Number of Pieces

Number of Pieces: Number of Pieces that you want to cut into. NetCDF file will save the time in its specific time step. There will be different format file. Thus, it is recommended to input pieces like 5 or 10

4.3.3 Number of Year

Number of Year: Time interval of the target file.

4.3.4 Number of time step

Number of time step: Number of time step. Needed because for different file, for same time interval, the time step might be different.

To see how many time-steps the target file have.

Run:

```
ncdump -h {target file path}
```

4.3.5 variable name

name of the variable you want to change unit

4.3.6 output unit

Output unit: new unit of the variable

4.3.7 Model Grid

The model grid that you want to re-grid the original into

4.3.8 output prefix

Output file's prefix to indicate the output

4.3.9 start year

Start year: get the start year of the target file to indicate the time range

4.3.10 output folder

Output folder: folder to store the file after preprocessing

4.3.10 regrid

Regrid : used to indicate whether to regrid or not

4.4 Input arguments for input file for Diagnostics

The following are the **required** arguments needed in the program input.

4.4.1 Start date of analysis

start_date : This can be in the following formats:

YYYY-MM-DD : e.g. 2020-04-12

YYYY-MM : e.g. 2020-04

YYYY : e.g. 2020

- If day is not given, the 1st of the given month will be used i.e. 2020-04 => 2020-04-01
- If day and month is not given, 1st Jan will be used as the start date i.e. 2020 => 2020-01-01

start_date1, start_date2 : 2 start dates can be put for control and future runs respectively. This is only used if [Calculate index](#) argument is set or [multi-model analysis between future and historical data](#) is performed .

In command line:

start_date or *start_date1 start_date2*

4.4.2 Prefix

prefix_of_file : This is the prefix of the file name.

prefix_of_file1, prefix_of_file2 : 2 prefixes can be put for control and future runs respectively. This is only used if [Calculate index](#) argument is set or [multi-model analysis between future and historical data](#) is performed.

If 2 prefixes are given, 2 [start dates](#) and 2 [end dates](#) **must** be set.

In command line:

`-pf prefix_of_file` or `-pf prefix_of_file1 prefix_of_file2`

For example, for file : `air_temp_1970.nc`, the prefix chosen can be `air` or `air_temp` or even `ai`. As long as the file names stored start with the prefix, it will be selected. You should ensure that if multiple model outputs are stored in `DATA/`, then the (unique) prefix of the files wanted should be used.

4.4.3 Variables

`var1, var2, ...` : Variables of data to analyse

In command line:

`-v var1 var2 ...`

The variables given must match the variable names in the files, otherwise an error will be thrown.

4.4.4 Number of ensembles

`num_of_ens` : The number of ensembles to analyse

In command line:

`-e num_of_ens`

The following are [optional](#) arguments that can be set as pleased. If any of these arguments are left blank, then it is automatically seen as false / not set.

4.4.5 End date of analysis

.

end_date : This is in the same format as the *start date*

If end date is not given:

- If only start year is given, the end date is automatically set to the 31 Dec of start year
- If start year and month is given, then end date is set to the end of the start month

If end date given:

- If day is not given, the end of the given month will be used i.e. 2020-04 => 2020-04-30
- If day and month is not given, 31 Dec will be used as the end date i.e. 2020 => 2020-12-31

end_date1, *end_date2*: **Must** be set if 2 *start dates* are given.

In command line:

`-end end_date or -end end_date1, end_date2`

4.4.6 Analysis

a1, *a2*, ... : Analysis performed on data

The types of analysis that can be selected:

- mean
- std (Standard deviation)
- rms (Root mean squared error)
- median

All the analysis calculated with respect to time, unless the *Spatial argument* is set.

In command line:

`-a a1 a2 ...`

Multi-model analysis between historical and future runs: If 2 start dates and 2 end dates are given, then the difference between the analysis between the time periods are computed. **Important:** The time steps of historical and future **must** be equal in order to successfully compute the difference, if analysis is set (without *spatial*). E.g. if you have start dates : 1850, 2010 and end dates : 1900, 2015, the program **will** fail with a Value error since the time period of the first date is 50, which is not the same as the time

period of the next date, 5. However, start dates : 1850, 2010 and end dates : 1855, 2015 will not fail as the time period is 5 years for both date sets.

4.4.7 Spatial

True/False : **Analysis** will be calculated spatially instead of temporally.

In command line:

-sp

4.4.8 Total ensemble stats

This setting is only applicable if multiple ensembles are given.

True/False : All ensembles are averaged (using mean) into one ensemble.

In command line:

-t

4.4.9 Plot

ens_num : Plot map, histogram and time series (and box plots) graphs given specific ensemble to plot. The ensemble number must be > 0. The plots show in pop up figures when program is run.

In command line:

-p *ens_num*

If only 1 ensemble given, then *ens_num* = 1.

If **Total ensemble stats** is set, then *ens_num* = 1 plots for the averaged ensemble.

When setting plot, the data given to plotting function will be the data passed through the **Analysis** argument. If the analysis argument is not set, then it will be the data with no analysis.

So if the data given to the plotting function is 2D (lat, lon), then the map is plotted. If data is 1D (time) then the time series and box plots are plotted. The histogram data will be plotted regardless of dimension of data (except some error is thrown).

4.4.10 Monthly

True/False : Set to True if data in file is stored in monthly increments, otherwise daily increments is assumed.

In command line:

-m

If there is only 1 time step in the data file, then setting Monthly has no effect.

4.4.11 Grid

latitude, longitude : Extracts the grid point that latitude and longitude lies in.

sample_points.txt : If a list of points are defined in a .txt file, then extracts the list of points. An example of what the file should contain is stored in INPUT/sample_points.txt.

example_file.nc : Regrids to the rectangular grid in NetCDF file given. This file should be stored in the DATA/ folder.

The interpolation scheme used is nearest neighbour.

In command line:

-g *latitude, longitude* or -g *sample_points.txt* or -g *example_file.nc*

Regridding only works between rectangular grids (this holds for [Sample points](#) as well). This setting **cannot** be used in conjunction with the [Sample](#) setting.

4.4.12 Sample

latitude, longitude : Extracts the sample point that latitude and longitude lies in.

sample_points.txt : If a list of points are defined in a .txt file, then extracts the list of points. An example of what the file should contain is stored in INPUT/sample_points.txt.

example_file.nc : Regrids to the rectangular grid in NetCDF file given. This file should be stored in the DATA/ folder.

The interpolation scheme used is linear.

In command line:

-s *latitude, longitude* or -s *sample_points.txt* or -s *example_file.nc*

Regridding only works between rectangular grids (this holds for [Grid points](#) as well). This setting **cannot** be used in conjunction with the [Grid](#) setting.

4.4.13 Mask file

mask_example.txt :Extracts region from data using a file that contains the description of multiple polygons. An example file is given in INPUT/mask_example.txt.

In command line:

```
-mk mask_example.txt
```

Example file in INPUT/mask_example.txt:

```
# Mask file expected to have only one list of polygons
# Comments in file can have '#' at the start of the line
# Each polygon is a list of tuples of integers (longitude, latitude)
# level is the depth level of mask. If this is not specified,
# then all levels of data are used.
# Note: level >= 1 and <= depth in NetCDF file
# You can also put a range of levels to choose separated by '-'
[(-25, -40), (75, -40), (-25, 40), (75,40)] level: 2
```

If data is 3D (time, lat, lon), level would select within the time dimensions. If the data is 4D (depth, time, lat, lon), then level selects within the depth dimension.

4.4.14 Save output

True/False : Saves data output in NetCDF files and saves tables of histogram and time series as .txt files. If *Plot* set, then saves the figures as .png as well.

In command line:

```
-O
```

4.4.15 Covary

True/False : Histogram co-variance. There must be only 2 variables in *Variables* argument. If *Plot* or *Save output* is set, then a 2D histogram plot is created.

In command line:

```
-CV
```

4.4.16 Histogram bin selection

hist_choice : Choose the number of bins used in the histogram. The default = fd, the Freedman Diaconis Estimator. To see a list of other estimators, go to this [the NumPy page](#).

num_of_bins: You can also put the number of bins e.g. 10, as the input.

num_of_bins1, num_of_bins2: If 2D histogram is being plotted, then the number of bins for each variable **must** be given.

histogram_bins.txt: If you want to specify the bin edges used in the histogram, you can input the file - an example file is given in INPUT/histogram_bins.txt

In command line:

```
-ht hist_choice or -ht num_of_bins or -ht num_of_bins1 num_of_bins2
```

4.4.17 Longitude centre

centre : Longitude to centre data on

In command line:

```
-lc centre
```

4.4.18 User function

file_name, *function_name* : Use function written by the user and stored in INPUT/user_function folder for analysis. The *file_name* is the name of file that contains function in user_function folder. The *function_name* is the name of function to call. There is an example file with functions in INPUT/user_function.

In command line:

```
-u file_name function_name
```

When the user function setting is turned on, then the [analysis](#) argument is automatically not used. If [Plot](#) argument is set, the graphs plotted are using only the original data (without the user function applied).

The user function argument does not consider the setting of the [Total ensemble stats](#) argument. This is because, the user function is applied to all ensembles.

Example file in INPUT/user_function/example_function.py

```
import iris
```

```
"""
```

This is a script that given an example of a user-given function in the command line or input file :

```
file_name = example_function
```

```
func_name = simple_equation
```

```
"""
```

```
def simple_equation(cube):
```

```
    """
```

Perform simple arithmetic with variables

:param data: a dictionary of variables and their data (iris.cube)

:return: result of equation, name of results, unit

```
    """
```

Variables are stored in the cube dictionary The names of variables used have to match ones in the Variables argument provided in the input file or command line.

```
temperature = cube['temp'].data
```

```
salinity = cube['sal'].data
```

```
# Perform equation with temperature and salinity result
```

```
= 2 * temperature + 4 * salinity
```

Set the attributes of data

```
name = 'result'
```

```
long_name = 'result calculated using simple equation' unit
```

```
= 'K'
```

```
return result, name, long_name, unit
```

4.4.19 Calculate areas

True/False : Calculate areas of grid boxes of latitude and longitude and saves to NetCDF file RESULTS/areas.nc.

In command line:

-ca

4.4.20 Calculate index

index : Calculate the index given The control run is the FIRST file *prefix* set and the corresponding *start/end* date. The future run is the SECOND file *prefix* set and the corresponding second *start/end* date. Types of indices that can be calculated:

- enso : The Oceanic Niño Index (ONI)
- nino12 : Niño 1+2 Index
- nino4 : Niño 4 Index
- tni : The Trans-Niño Index (TNI)
- iod : Indian Ocean Dipole (IOD) Mode Index
- amo : Atlantic Multidecadal Oscillation (AMO) Index
- pdo : Pacific Decadal Oscillation (PDO) Index
- ao : Arctic Oscillation (AO; Northern Annular Mode) Index
- aao : Antarctic Oscillation (AAO; Southern Annular Mode) Index
- nao : North Atlantic Oscillation (NAO) Index

In command line:

-i *index*

Important: One file for control and one file for future run is expected.

If multiple files are used for the control run, then combine them into one first, this *application* gives the input arguments necessary.

In order to use the Calculate index functionality, *CDO* needs to be installed, since the indices are calculated using this package.

1. Go to <https://code.mpimet.mpg.de/projects/cdo/files>
2. Download the most recent **cdo-1.*.*.tar.gz** file

3. On the command line, go to the place where it is downloaded.

4. Extract data from tar zip file:

```
$ tar xvfz cdo -1 . * . * . * . tar.gz
```

5. cd into newly extracted directory

```
$ cd cdo -1 . * . * . * . /
```

6. do:

```
$ make
```

7. do:

```
$ make install
```

8. To check if installation has happened successfully, do

```
$ cdo -V
```



Some applications

The following are some scenarios.

5.0.1 Extract sea surface temperature between 1980 and 1990 at grid point (14.5, -27.5) and output a time series, histogram and NetCDF file with mean data

REQUIRED ARGUMENTS

Prefix: sst

Start date of analysis: 1980

Variables: sst

Number of ensembles: 1

#

OPTIONAL ARGUMENTS

End date of analysis: 1990

Analysis: mean

Spatial:

Total ensemble stats:

Plot: 1

Monthly:

Grid: 14.5, -27.5

Sample:

Mask file:

Save Output: True

Covary:

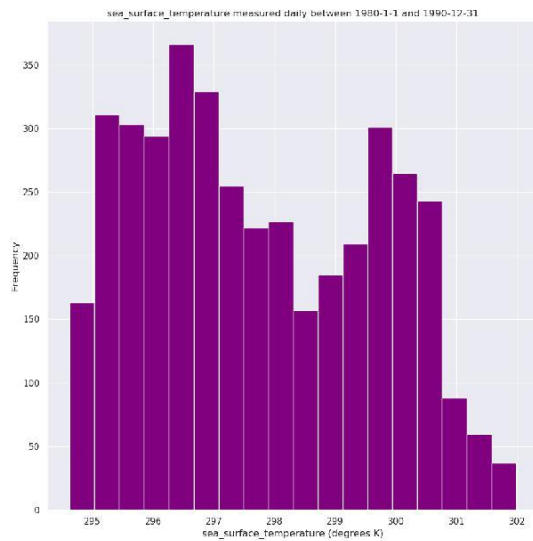
Histogram bin selection:

Longitude centre:

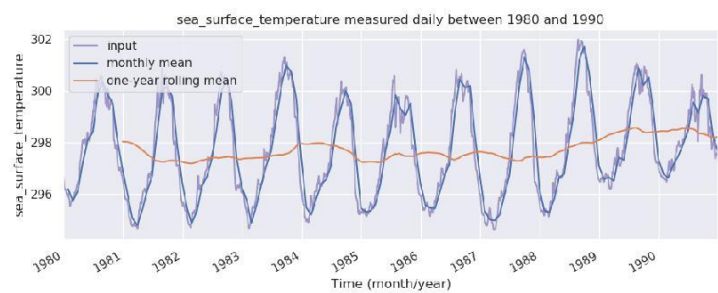
User function:

Calculate areas:

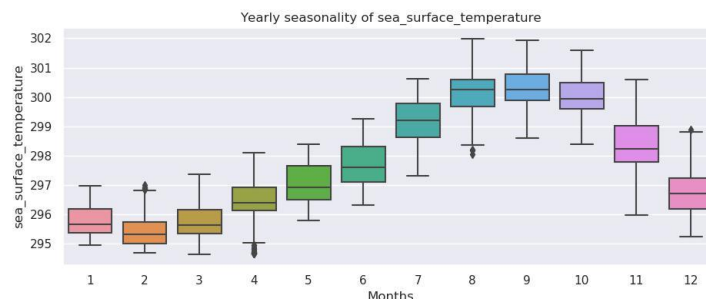
Calculate index:



(a) Histogram



(b) Time series

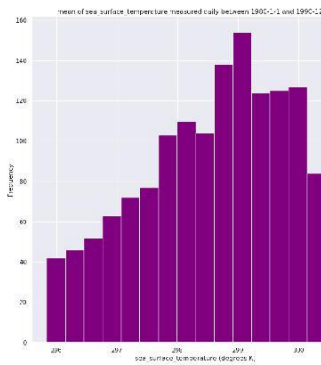


(c) Box plots

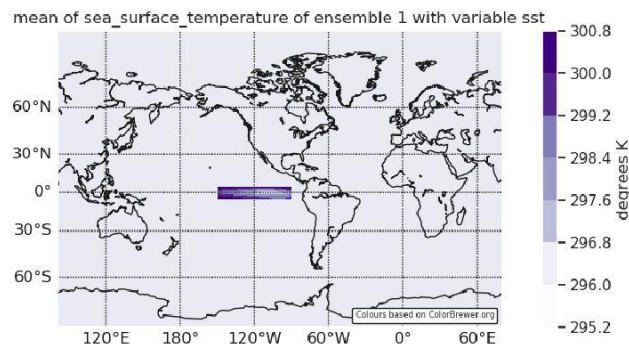
Figure 1: Plots generated for scenario 5.0.1

5.0.2 Calculate mean sea surface temperature between 1980 and 1990 using region of NINO-3 and output histogram, map and NetCDF file with mean data

```
# REQUIRED ARGUMENTS
# -----
Prefix: sst
Start date of analysis: 1980
Variables: sst
Number of ensembles: 1
#
# -----
# OPTIONAL ARGUMENTS
# -----
End date of analysis: 1990
Analysis: mean
Spatial:
Total ensemble stats:
Plot: 1
Monthly:
Grid:
Sample:
Mask file: nino3_mask.txt
Save Output: True
Covary:
Histogram bin selection:
Longitude centre:
User function:
Calculate areas:
Calculate index:
```



(a) Histogram

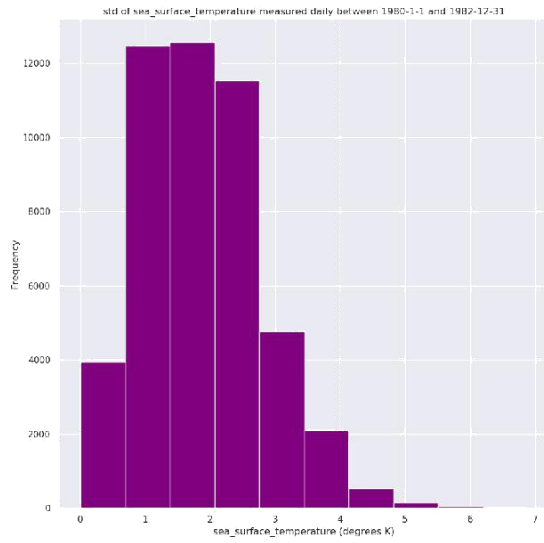


(b) Map with NINO-3

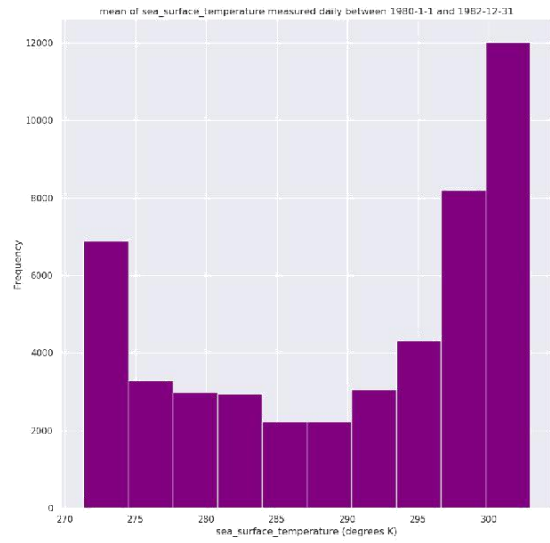
Figure 2: Plots generated for scenario 5.0.2

5.0.3 Calculate mean and std of sea surface temperature between 1980 and 1982, with longitude centred at -30 degrees east. Histogram bins = 10 and grid boxes areas calculated. Outputs are NetCDF files containing areas and mean and std.

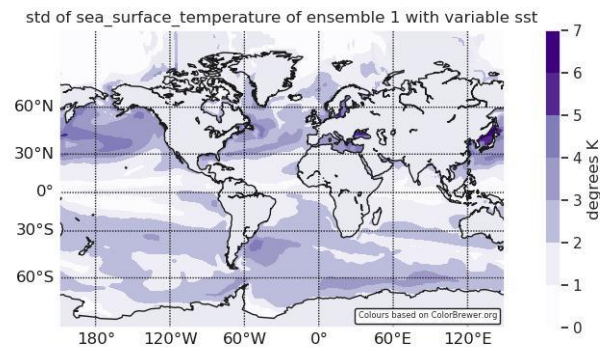
```
# REQUIRED ARGUMENTS
# -----
Prefix: sst
Start date of analysis: 1980
Variables: sst
Number of ensembles: 1
#
# -----
# OPTIONAL ARGUMENTS
# -----
End date of analysis: 1982
Analysis: mean, std
Spatial:
Total ensemble stats:
Plot:
Monthly:
Grid:
Sample:
Mask file:
Save Output:
Covary:
Histogram bin selection: 10
Longitude centre: -30
User function:
Calculate areas: True
Calculate index:
```

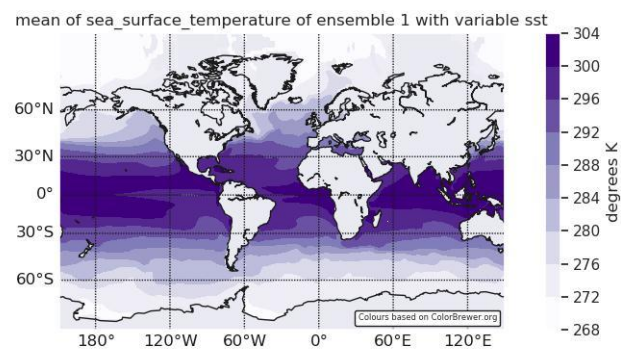
(a) Histogram of Standard deviation



(b) Histogram of mean



(c) Map of Standard deviation



(d) Map of mean

Figure 3: Plots generated

5.0.4 Combine consecutive files of sea surface temperature into one

Given data files

- sst_1950_1959.nc
- sst_1960_1969.nc
- sst_1970_1979.nc

REQUIRED ARGUMENTS

Prefix: sst

Start date of analysis: 1950

Variables: sst

Number of ensembles: 1

#

OPTIONAL ARGUMENTS

End date of analysis: 1979

Analysis:

Spatial:

Total ensemble stats:

Plot:

Monthly:

Grid:

Sample:

Mask file:

Save Output: True

Covary:

Histogram bin selection:

Longitude centre:

User function:

Calculate areas:

Calculate index:

Output file generated: sst_1950_1979.nc.

For more applications/scenarios, go to the [integration tests section](#).



Tests

6.1 Unit tests

The testing framework [Pytest](#) was used to construct the unit tests written. In order to run the unit tests, Pytest has to be installed with:

```
$ pip install pytest
```

After installation,

```
$ cd EZclim/
```

and finally run:

```
$ pytest
```

You should get output similar to:

```
(base) ada@DESKTOP-9T9CN93:/mnt/c/Users/Adanna/acse-9-independent-research-project-AdannaAkwats/ClimateModelling$ pytest
===== test session starts =====
platform linux -- Python 3.7.3, pytest-3.5.1, py-1.8.0, pluggy-0.6.0
rootdir: /mnt/c/Users/Adanna Akwataghibe/acse-9-independent-research-project-AdannaAkwats/ClimateModelling, inifile:
collected 33 items

tests/test_user_entry.py ..... [ 18%]
tests/test_utils.py ..... [100%]

===== 33 passed in 2.17 seconds =====
```

Figure 4: Pytest command line output

6.2 Integration tests

We did some *scenario* testing, where we constructed test cases, ran them and generated an expected output. During the development of the code, these scenarios were run frequently to ensure that the actual output matched the expected output. The tests are stored in EZclim/integration_tests.

Integration tests folder structure

```
EZclim/integration_tests/  
|-- Test1_mean_spatial_monthly/  
|-- Test2_std_monthly/  
|   |-- data/  
|   |-- results/  
|   |-- plots/  
|   |-- input.txt  
|   |-- other input files ...  
|   ...
```

- data/ contains the NetCDF data files used when running the program
- results/ contains the programs NetCDF outputs
- plots/ contains the .png and .txt files for maps/histogram/time series generated.
- the input files consists of the input.txt file as well as any mask.txt, sample.txt, histogram.txt files used.

You can use the data and the input given to run some applications. And then check that the output generated matches the plots and results.

The test cases

1. Test_mean_spatial_monthly : Calculate the spatial mean of air pressure at convective cloud base between 1979 and 2014 measured monthly, with histogram bin size of 10.
2. Test_std_monthly : Calculate the standard deviation of air pressure at convective cloud base between 1979 and 2014 measured monthly.
3. Test_multimodel_median_monthly : Calculate the multi model median difference for air pressure at convective cloud base measured monthly, between 1950-1959 (control) and 2010-2014.
4. Test_multimodel_mean_spatial_monthly : Calculate the multi model spatial mean difference of air pressure at convective cloud base measured monthly, between 1950-1959 (control) and 2010-2014.

5. `Test_enso_index_monthly` : Calculate the ENSO index for air pressure at convective cloud base measured monthly, between 1950-1959 (control) and 2010-2014.
6. `Test_gridpoint_monthly_areas`: Extract the grid point (20, 20) for air pressure at convective cloud base between 1979 and 2014 measured monthly.
7. `Test_samplefile`: Regrid air pressure at convective cloud base between 1979 and 2014 measured monthly to grid of model output for North America.
8. `Test_mask_mean_monthly_centred_0.0` : Calculate the mean of air pressure at convective cloud base between 1979 and 2014 measured monthly within region given by the mask and with the longitude centred at 0.0.
9. `Test_mean_total_spatial` Calculate the spatial mean of 2 ensembles of air temperature at 1970, giving one average over all ensembles.
10. `Test_mean_4D_mask_2vars` : Calculate the mean of temperature and relative humidity at 2193 within the region given by mask. (The grid used is 4D (depth, time, lat, lon)).
11. `Test_4D_user_function` : Perform equation between temperature and relative humidity given by user function of temperature and relative humidity at 2193. (The grid used is 4D (depth, time, lat, lon)).
12. `Test_4D_2DHist` : Calculate the mean of temperature and relative humidity and create 2D histogram between the variables.
13. `Test_4D_2DHist_file` : Create 2D histogram between temperature and relative humidity using user defined histogram bins.
14. `Test_total` : Combine 2 ensembles of air temperature at 1970, giving one average over all ensembles.

Running on different setups (running locally)

The program was set up and run on different types of machines:

1. PC
2. Workstation (ssh-ed into Linux workstation)
3. HPC (High Power Computing) cluster

If you are running the program on a PC or Workstation, then the program should run without any additional setup.

Running in a HPC requires a little more care. The parallelised version of the code is stored in the EZclim_Parallel/ folder. There are some requirements in order to run on HPC:

7.0.1 Storage

When using multiple ensembles (especially for already large model outputs), the output given by the program can get very big - depending on the analysis being done. In some HPC systems, the storage used for the cluster job needs to be specified ahead of time. The size of the input files + the output needs to be taken into account when specifying the space allocated for the job.

7.0.2 Number of Processors

The number of processors used in the code, can be set in the file EZclim_Parallel/parallel_setting.py.

7.0.3 Plotting

On a NPC cluster, the graphical display for plots is not available. Therefore, the [Plotting argument](#) **must not** be set, in order for the program to run successfully.

For example, [this application](#) would need to be changed to the input file below to work on a HPC.

```
# REQUIRED ARGUMENTS
# -----
Prefix: sst
Start date of analysis: 1980
Variables: sst
Number of ensembles: 1
#
# -----
# OPTIONAL ARGUMENTS
# -----
End date of analysis: 1990
Analysis: mean
Spatial:
Total ensemble stats:
Plot:          # The plot is not set
Monthly:
Grid:
Sample:
Mask file: nino3_mask.txt
Save Output: True
Covary:
Histogram bin selection: fd
Longitude centre:
User function:
Calculate areas:
Calculate index:
```




Troubleshooting

Below are some bugs found in the code that you may encounter.

8.0.1 (Seemingly) Non-unique prefix

Say you have 2 files `air_temp_1970.nc` and `air_temp2_1970.nc` and only wanted to analyse the first one. Setting the `Prefix` argument to `air_temp` selects both files instead of just the files. Instead just use the prefix: `air_temp_`. You just need to choose a unique prefix to the first file. The entire file name can be chosen as a prefix if you are in a special case where the file names are too similar.

8.0.2 Number of histogram bins not equal to what selected

When choosing the number of bins as an integer, the number may not be exactly the same as what you set as the argument, the `NumPy function` used to plot histogram seems to do some sort of auto bin calculation.

Alternatively, if one of the estimators is chosen, sometimes the number of bins generated causes a memory error. If this is the case, the program reverts to plotting the histogram with about 10 bins.

8.0.3 Changes to dimension names

When saving the output file, the Python package `xarray` sometimes saves dimensions as `dim_0` and `dim_1`, not by the original dims names longitude/latitude.

8.0.4 Plot not displayed on Windows 10

In order to display plots in Windows (from the Ubuntu subsystem), you need to download Xming for Windows from <https://sourceforge.net/projects/xming/>. Then

1. Start XLaunch
2. Select **Multiple Windows**
3. Set **Display number** to **0**
4. Click Next
5. Select **Start no client** option
6. Click Next
7. Do not change anything in **Specify Parameter settings** , Click Next again.
8. Save configuration - Click on this saved file to start set up XLaunch if you log out of your system.

Now that the connection is made, go to the linux cmd and type

```
$ export DISPLAY= : 0
```

Now running the code with plotting on should give a display.

If you are running on a Linux terminal using ssh, then this [web-page linked \(http://laptops.eng.uci.edu/software-installation/using-linux/how-to-configure-xming-putty\)](http://laptops.eng.uci.edu/software-installation/using-linux/how-to-configure-xming-putty) gives the steps to connect to Xming when ssh-ing using putty.

8.0.5 Permission error given when writing/finding file

If a permission error is thrown when writing/finding file, make sure there are no spaces within your file path.

8.0.6 ValueError: Variable '..' has multiple fill values

This is a problem with the Python package [xarray](#). You will need to remove the 'missing value' attribute since xarray fails when both fill and missing value are set.

8.0.7 Processes when running in parallel unable to transfer data

The code is parallelised between ensemble data. So each processor does work for a single ensemble. This means that if the data is too big, then transferring data between processes can become an issue. This can lead to the code throwing a `multiprocess.pool memory error`, or the code may stay in a hanged state - and take a *very* long time to load.

8.0.8 AttributeError: attributes of masked are not writeable

If this error is thrown: *AttributeError: attributes of masked are not writeable*, then the sample/grid point or mask has values of all NaNs. A new point or mask regions that is not encompassed just of NaNs needs to be chosen.

8.0.9 ValueError: operands could not be broadcast together with shapes (,) (**)**

If you are trying to perform multi-model analysis, the error is most likely due to the different time period between the control and future run. Go to the [Multi model section](#) for more help.