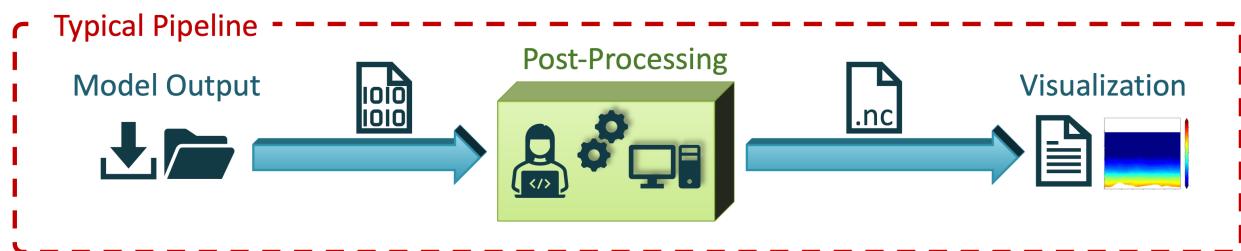
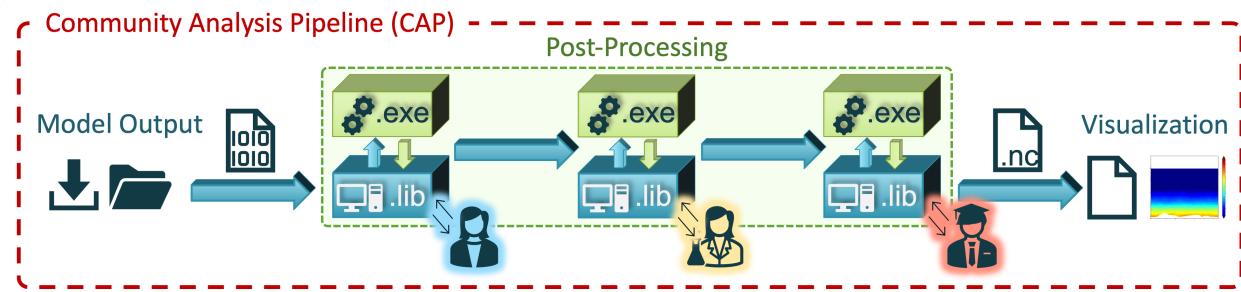


## Introducing the Community Analysis Pipeline (CAP)

CAP is a toolkit designed to simplify the post-processing of MGCM output. CAP is written in Python and works with existing Python libraries, allowing any Python user to install and use CAP easily and free of charge. Without CAP, plotting MGCM output requires that a user provide their own scripts for post-processing, including code for interpolating the vertical grid, computing and adding derived variables to files, converting between file types, and creating diagnostic plots. In other words, a user would be responsible for the entire post-processing effort as illustrated in Figure 1.



Such a process requires that users be familiar with Fortran files and be able to write (or provide) script(s) to perform file manipulations and create plots. CAP standardizes the post-processing effort by providing executables that can perform file manipulations and create diagnostic plots from the command line. This enables users of almost any skill level to post-process and plot MGCM data (Figure 2).



As a foreword, we will list a few design characteristics of CAP:

- CAP is written in **Python**, an open-source programming language with extensive scientific libraries available
- CAP is installed within a Python **virtual environment**, which provides cross-platform support (MacOS, Linux and Windows), robust version control (packages updated within the main Python distribution will not affect CAP), and is not intrusive as it disappears when deactivated
- CAP is composed of a **set of libraries** (functions), callable from a user's own scripts and a collection of **five executables**, which allows for efficient processing of model outputs from the command-line.
- CAP uses the **netCDF4 data format**, which is widely used in the climate modeling community and self-descriptive (meaning that a file contains explicit information about its content in terms of variable names, units etc...)
- CAP uses a convention for output formatting inherited from the GFDL Finite-Volume Cubed-Sphere Dynamical Core, referred here as "**FV3 format**": outputs may be binned and averaged in time in various ways for analysis.
- CAP long-term goal is to offer **multi-model support**. At the time of writing, both the NASA Ames Legacy GCM and the NASA Ames GCM with the FV3 dynamical core are supported. Efforts are underway to offer compatibility to other Global Climate Models (e.g. eMARS, LMD, MarsWRF).

Specifically, CAP consists of five executables:

1. MarsPull.py Access MGCM output
2. MarsFiles.py Reduce the files
3. MarsVars.py Perform variable operations
4. MarsInterp.py Interpolate the vertical grid
5. MarsPlot.py Visualize the MGCM output

These executables and their commonly-used functions are illustrated in the cheat sheet below in the order in which they are most often used. You should feel free to reference during and after the tutorial.

## Cheat sheet

---



NASA AMES

MARS CLIMATE MODELING CENTER

## Quick Start

### Create, Source the Environment:

```
% /path/to/python3 -m venv -system-site-packages amesGCM3  
% source ~/amesGCM3/bin/activate
```

### Install CAP:

```
% pip install git+https://github.com/alex-kling/amesgcm.git
```

### Update CAP:

```
% pip install git+https://github.com/alex-kling/amesgcm.git --upgrade
```

### Uninstall CAP:

```
% pip uninstall amesgcm
```

### Deactivate the Environment:

```
% deactivate
```

## Frequently Used Commands



### Access MGCM Output

```
% MarsPull.py -id INERTCLDS -ls 255 285  
% MarsPull.py -id INERTCLDS -f fort.11_0670 fort.11_0671
```



### File Manipulations

```
% MarsFiles.py LegacyGCM_*.nc -fv3  
% MarsFiles.py *atmos_average.nc -combine  
% MarsFiles.py *atmos_diurn.nc -tshift
```



### Variable Operations

```
% MarsVars.py *atmos_average.nc -col vap_mass  
% MarsVars.py *atmos_average.nc -add rho
```



### Interpolations

```
% MarsInterp.py *atmos_average.nc -t pstd  
% MarsInterp.py *atmos_average.nc -t pstd -l phalf_mb
```



### Visualizations and File Contents

```
% MarsPlot.py -template  
% MarsPlot.py Custom.in  
% MarsPlot.py -i 00000.atmos_average.nc
```

CAP is designed to be modular. For example, a user could post-process and plot MGCM output exclusively with CAP or a user could employ their own post-processing routine and then use CAP to plot the data. Users are free to selectively integrate CAP into their own analysis routine to the extent they see fit.

## Table of Contents

- Introducing the Community Analysis Pipeline (CAP)

- Cheat sheet
- The big question... How do I do this? > **Ask for help!**
- 1. MarsPull.py - Downloading Raw MGCM Output
- 2. MarsFiles.py - Reducing the Files
- 3. MarsVars.py - Performing Variable Operations
- 4. MarsInterp.py - Interpolating the Vertical Grid
- 5. MarsPlot.py - Plotting the Results
- MarsPlot.py: How to?
  - Disable or add a new plot
  - Adjust the color range and colormap
  - Make a 1D-plot
  - Customize 1D plots
  - Put multiple plots on the same page
  - Put multiple 1D-plots on the same page
  - Use a different epoch
  - Access simulation in a different directory
  - Overwrite the free dimensions.
  - Element-wise operations
  - Commenting out and speed-up processing
  - Change projections
  - Figure format, size
  - Access CAP libraries and make your own plots
  - Debugging

---

## The big question... How do I do this? > **Ask for help!**

---

Use the `--help` (`-h` for short) option on any executable to display documentation and examples.

```
(amesGCM3)>$ MarsPlot.py -h
> usage: MarsPlot.py [-h] [-i INSPECT_FILE] [-d DATE [DATE ...]] [--template]
>                      [-do DO] [-sy] [-o {pdf,eps,png}] [-vert] [-dir DIRECTORY]
>                      [--debug]
>                      [custom_file]
```

---

## 1. MarsPull.py - Downloading Raw MGCM Output

---

MarsPull is a utility for accessing MGCM output files hosted on the [MCMC Data portal](#). MGCM data is archived in 1.5 hour intervals (16x/day) and packaged in files containing 10 sols. The files are named fort.11\_XXXX in the order they were produced, but MarsPull maps those files to specific solar longitudes ( $L_s$ , in  $^{\circ}$ ). This allows users to request a file at a specific  $L_s$  or for a range of  $L_s$  using the `-ls` flag. Additionally the `identifier` (`-id`) flag is used to route MarsPull through a particular simulation. The `filename` (`-f`) flag can be used to parse specific files within a particular directory.

```
MarsPull.py -id INERTCLDS -ls 255 285
MarsPull.py -id ACTIVECLDS -f fort.11_0720 fort.11_0723
```

[Back to Top](#)

---

## 2. MarsFiles.py - Reducing the Files

MarsFiles provides several tools for file manipulations, including code designed to create binned, averaged, and time-shifted files from MGCM output. The `-fv3` flag is used to convert fort.11 binaries to the Netcdf data format (you can select one or more of the file format listed below):

```
(amesGCM3)>$ MarsFiles.py fort.11* -fv3 fixed average daily diurn
```

These are the file formats that MarsFiles can create from the fort.11 MGCM output files.

### Primary files

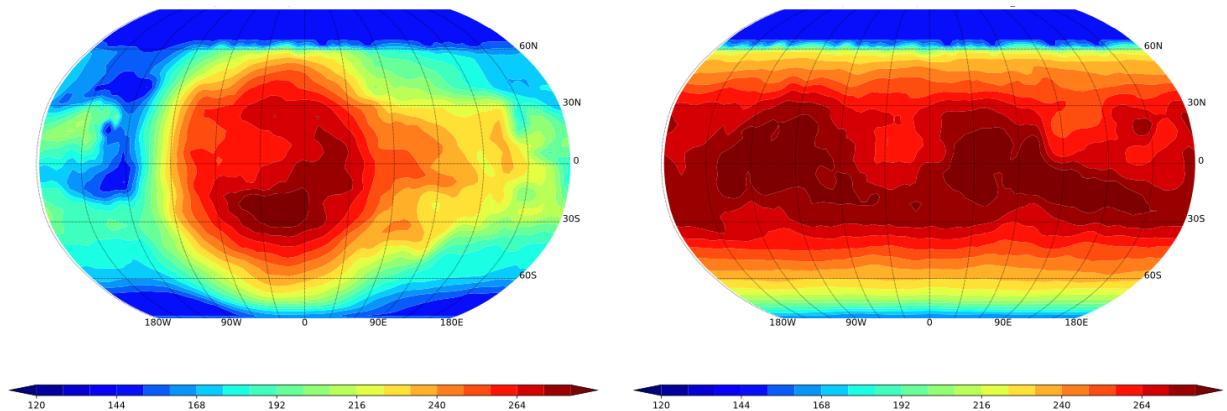
| File name               | Description                                            | Timesteps for 10 sols x 16 output/sol | Ratio to daily file (430Mb) |
|-------------------------|--------------------------------------------------------|---------------------------------------|-----------------------------|
| <b>atmos_daily.nc</b>   | continuous time series                                 | (16 x 10)=160                         | 1                           |
| <b>atmos_diurn.nc</b>   | data binned by time of day and 5-day averaged          | (16 x 2)=32                           | x5 smaller                  |
| <b>atmos_average.nc</b> | 5-day averages                                         | (1 x 2) = 2                           | x80 smaller                 |
| <b>fixed.nc</b>         | statics variable such as surface albedo and topography | static                                | few kB                      |

### Secondary files

| File name                        | description                                             |
|----------------------------------|---------------------------------------------------------|
| <b>daily_lpf,_hpfp,_bpfp</b>     | low, high and band pass filtered                        |
| <b>diurn_T</b>                   | uniform local time (same time of day at all longitudes) |
| <b>diurn_tidal</b>               | tidally-decomposed files into harmonics                 |
| <b>daily_to_average_to_diurn</b> | custom re-binning of daily files                        |

- MarsFiles can concatenate like-files together on the time dimension using the `-combine` (`-c`) flag.

```
> 07180.atmos_average.nc 07190.atmos_average.nc 07200.atmos_average.nc # 3 files with 10 days of output each  
(amesGCM3)>$ MarsFiles.py *atmos_average.nc -c  
> 07180.atmos_average.nc # 1 file with 30 days of output
```



3pm surface temperature before (left) and after (right) processing a diurn file with MarsFile to uniform local time

[Back to Top](#)

### 3. MarsVars.py - Performing Variable Operations

`MarsVars` provides several tools relating to variable operations such as adding and removing variables and performing column integrations. With no other arguments, passing a file to `MarsVars` displays file content much like `ncdump`:

```
(amesGCM3)>$ MarsVars.py 00000.atmos_average.nc
>
> =====DIMENSIONS=====
> ['bnds', 'time', 'lat', 'lon', 'pfull', 'scalar_axis', 'phalf']
> (etc)
> =====CONTENT=====
> pfull      : ('pfull',)= (30,), ref full pressure level [Pa]
> temp       : ('time', 'pfull', 'lat', 'lon')= (4, 30, 180, 360), temperature [K]
> (etc)
```

A typical option of `MarsVars` would be to add the atmospheric density `rho` to a file. Because the density is easily computed from the pressure and temperature fields, we do not archive it in the GCM output and instead provides a utility to add it as needed. This conservative approach to logging output allows to minimize disk space and speed-up post processing.

```
(amesGCM3)>$ MarsVars.py 00000.atmos_average.nc --add rho
```

We can see that `rho` was added by calling `MarsVars` with no argument as before:

```
(amesGCM3)>$ MarsVars.py 00000.atmos_average.nc
>
> =====DIMENSIONS=====
> ['bnds', 'time', 'lat', 'lon', 'pfull', 'scalar_axis', 'phalf']
> (etc)
> =====CONTENT=====
> pfull      : ('pfull',)= (30,), ref full pressure level [Pa]
> temp       : ('time', 'pfull', 'lat', 'lon')= (4, 30, 180, 360), temperature [K]
> rho        : ('time', 'pfull', 'lat', 'lon')= (4, 30, 180, 360), density (added postprocessing) [kg/m3]
```

The `help (-h)` option provides information on available variables and needed fields for each operation.

```
> Usage: MarsVars ****.atmos.average.nc -add rho
ON NATIVE FILES:
rho      (density)          Req. [ps,temp]
theta    (pot. temperature)  Req. [ps,temp]
pfull3D  (pressure at layer midpoint)  Req. [ps,temp]
DP       (layer pressure thickness)   Req. [ps,temp]
zfull    (altitude AGL)        Req. [ps,temp]
DZ       (layer altitude thickness)  Req. [ps,temp]
w        (vertical winds)       Req. [ps,temp,omega]
wdir     (wind direction)      Req. [ucomp,vcomp]
wspeed   (wind magnitude)     Req. [ucomp,vcomp]
N        (Brunt Vaisala freq)  Req. [ps,temp]
Ri       (Richardson number)   Req. [ps,temp]
Tco2     (CO2 condensation temperature)  Req. [ps,temp]
scorer_wl (Scorer horizontal wavelength)  Req. [ps,temp,ucomp]
div      (divergence)         Req. [ucomp,vcomp]
curl     (relative vorticity)   Req. [ucomp,vcomp]
fn       (frontogenesis)      Req. [ucomp,vcomp,theta]
```

#### NOTE:

Some support on interpolated files, in particular if pfull3D and zfull are added before interpolation to \_pstd, \_zagl, \_zstd.

#### ON INTERPOLATED FILES :

|      |                                    |                    |
|------|------------------------------------|--------------------|
| msf  | (mass stream function)             | Req. [vcomp]       |
| ep   | (wave potential energy)            | Req. [temp]        |
| ek   | (wave kinetic energy)              | Req. [ucomp,vcomp] |
| mx   | (vertical flux of zonal momentum)  | Req. [ucomp,w]     |
| my   | (vertical flux of merid. momentum) | Req. [vcomp,w]     |
| ax   | (zonal wave-mean flow forcing)     | Req. [ucomp,w,rho] |
| ay   | (merid. wave-mean flow forcing)    | Req. [ucomp,w,rho] |
| tp_t | (norm. temperature perturbation)   | Req. [temp]        |

MarsVars also offers the following variable operations:

| Command       | flag     | action                                                     |
|---------------|----------|------------------------------------------------------------|
| add           | -add     | add a variable to the file                                 |
| remove        | -rm      | remove a variable from a file                              |
| extract       | -extract | extract a list of variables to a new file                  |
| col           | -col     | column integration, applicable to mixing ratios in [kg/kg] |
| zdiff         | -zdiff   | vertical differentiation (e.g. compute gradients)          |
| zonal_detrend | -zd      | zonally detrend a variable                                 |

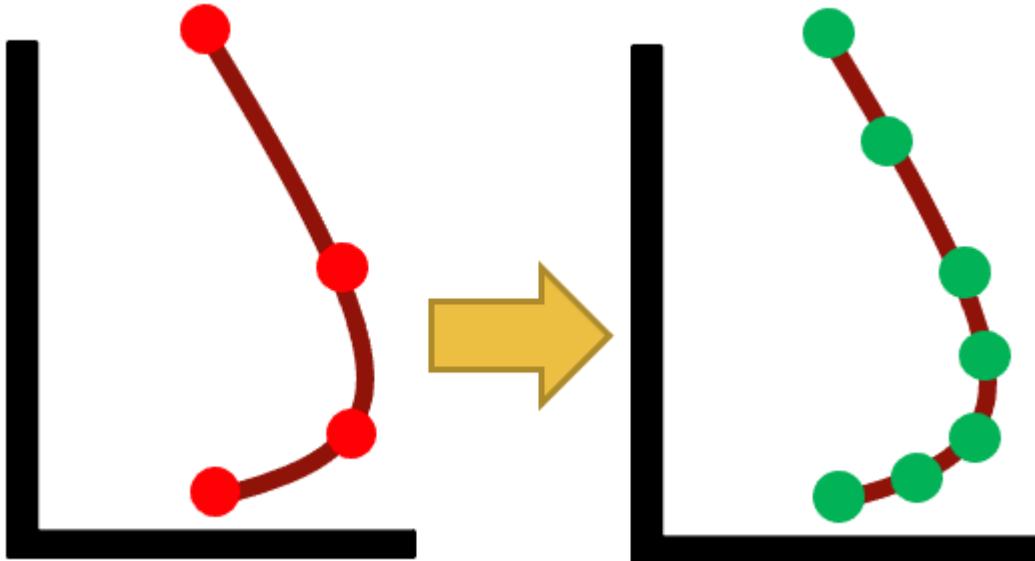
[Back to Top](#)

---

## 4. MarsInterp.py - Interpolating the Vertical Grid

---

Native MGCM output files use pressure as the vertical coordinate (`pfull`), which means the geometric height and pressure level of an atmospheric layer varies based on location.



#### *Pressure interpolation from the reference pressure grid to a standard pressure grid*

Climate data is usually analyzed on a standardized grid, however, and it is often necessary to interpolate the files to standard pressure coordinates. The `-type` (`-t`) argument in `MarsInterp` can interpolate files for you:

```
(amesGCM3)>$ MarsInterp.py 00000.atmos_average.nc -t pstd
```

An inspection of the file shows that the pressure level axis which was `pfull` (30 layers) has been replaced by a standard pressure coordinate `pstd` (36 layers), and all 3- and 4-dimensional variables reflect the new shape:

```
(amesGCM3)>$ MarsInterp.py 00000.atmos_average.nc -t pstd
(amesGCM3)>$ MarsVars.py 00000.atmos_average_pstd.nc
>
> =====DIMENSIONS=====
> ['bnds', 'time', 'lat', 'lon', 'scalar_axis', 'phalf', 'pstd']
> =====CONTENT=====
> pstd      : ('pstd',)= (36,), pressure [Pa]
> temp      : ('time', 'pstd', 'lat', 'lon')= (4, 36, 180, 360), temperature [K]
```

The following `type` (`-t`) flag of vertical interpolation are supported:

| file type          | description                              | low-level value in a deep crater |
|--------------------|------------------------------------------|----------------------------------|
| <code>_pstd</code> | standard pressure [Pa]                   | 1000Pa                           |
| <code>_zstd</code> | standard altitude [m]                    | -7000m                           |
| <code>_zagl</code> | standard altitude above ground level [m] | 0 m                              |

#### **Use of custom vertical grids**

It is also possible for the users to specify the layers for the interpolation. This is done by editing a **hidden** file `.amesgcm_profile` (note the dot ' . ') in your home directory.

For the first use, you will need to copy a template of `amesgcm_profile` to your /home directory:

```
(amesGCM3)>$ cp ~/amesGCM3/mars_templates/amesgcm_profile ~/.amesgcm_profile # Note the dot '.' !!!
```

You can open `~/.amesgcm_profile` with any text editor:

```
> <<<<<<<<| Pressure definitions for pstd |>>>>>>>>>
>p44=[1.0e+03, 9.5e+02, 9.0e+02, 8.5e+02, 8.0e+02, 7.5e+02, 7.0e+02,
>      6.5e+02, 6.0e+02, 5.5e+02, 5.0e+02, 4.5e+02, 4.0e+02, 3.5e+02,
>      3.0e+02, 2.5e+02, 2.0e+02, 1.5e+02, 1.0e+02, 7.0e+01, 5.0e+01,
>      3.0e+01, 2.0e+01, 1.0e+01, 7.0e+00, 5.0e+00, 3.0e+00, 2.0e+00,
>      1.0e+00, 5.0e-01, 3.0e-01, 2.0e-01, 1.0e-01, 5.0e-02, 3.0e-02,
>      1.0e-02, 5.0e-03, 3.0e-03, 5.0e-04, 3.0e-04, 1.0e-04, 5.0e-05,
>      3.0e-05, 1.0e-05]
>
>phalf_mb=[50]
```

In the example above, the user custom-defined two vertical grids, one with 44 levels (named `p44`) and one with a single layer at 50 Pa =0.5mbar(named `phalf_mb` )

You can use these by calling `MarsInterp` with the `-level` (`-l`) argument followed by the name of the new grid in `amesgcm_profile` .

```
(amesGCM3)>$ MarsInterp.py 00000.atmos_average.nc -t pstd -l p44
```

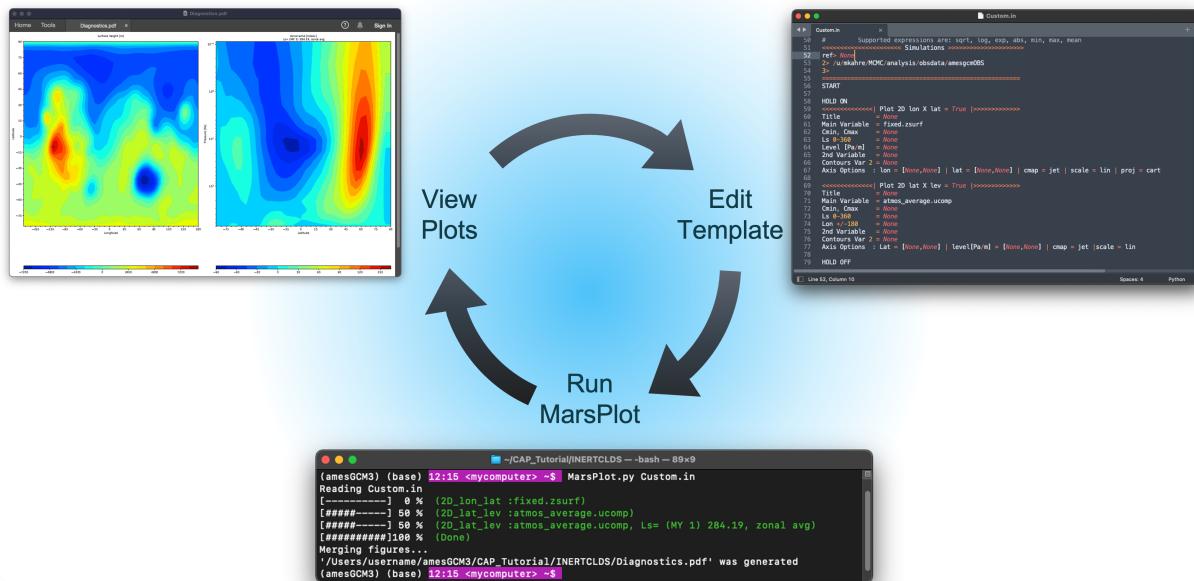
[Back to Top](#)

## 5. `MarsPlot.py` - Plotting the Results

The last component of CAP is the plotting routine, `MarsPlot` , which accepts a modifiable template ( `Custom.in` ) containing a list of plots to create. `MarsPlot` is useful for creating plots from MGCM output quickly, and it is designed specifically for use with the `netCDF` output files ( `daily` , `diurn` , `average` , `fixed` ).

The following figure shows the three components of `MarsPlot`:

- `MarsPlot.py`, opened in a **terminal** to inspect the netcdf files and ingest the `Custom.in` template
- `Custom.in` , a template opened in a **text editor**
- `Diagnostics.pdf`, refreshed in a **pdf viewer**



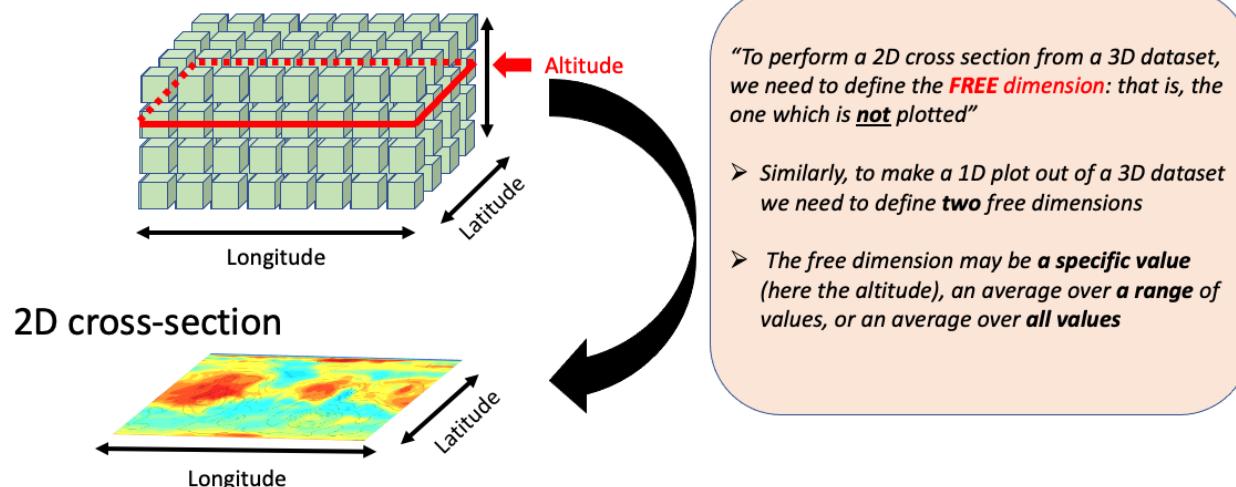
The default template, `Custom.in`, can be created by passing the `-template` argument to `MarsPlot`. `Custom.in` is pre-populated to draw two plots on one page: a topographical plot from the fixed file and a cross-section of the zonal wind from the average file. Creating the template and passing it into `MarsPlot` creates a PDF containing the plots:

```
(amesGCM3)>$ MarsPlot.py -template
> /path/to/simulation/run_name/history/Custom.in was created
(amesGCM3)>$
(amesGCM3)>$ MarsPlot.py Custom.in
> Reading Custom.in
> [-----] 0 % (2D_lon_lat :fixed.zsurf)
> [#####----] 50 % (2D_lat_lev :atmos_average.ucomp, Ls= (MY 2) 252.30, zonal avg)
> [#####----]100 % (Done)
> Merging figures...
> /path/to/simulation/run_name/history/Diagnostics.pdf was generated
```

Specifically MarsPlot is designed to generate 2D cross - sections and 1D plots.

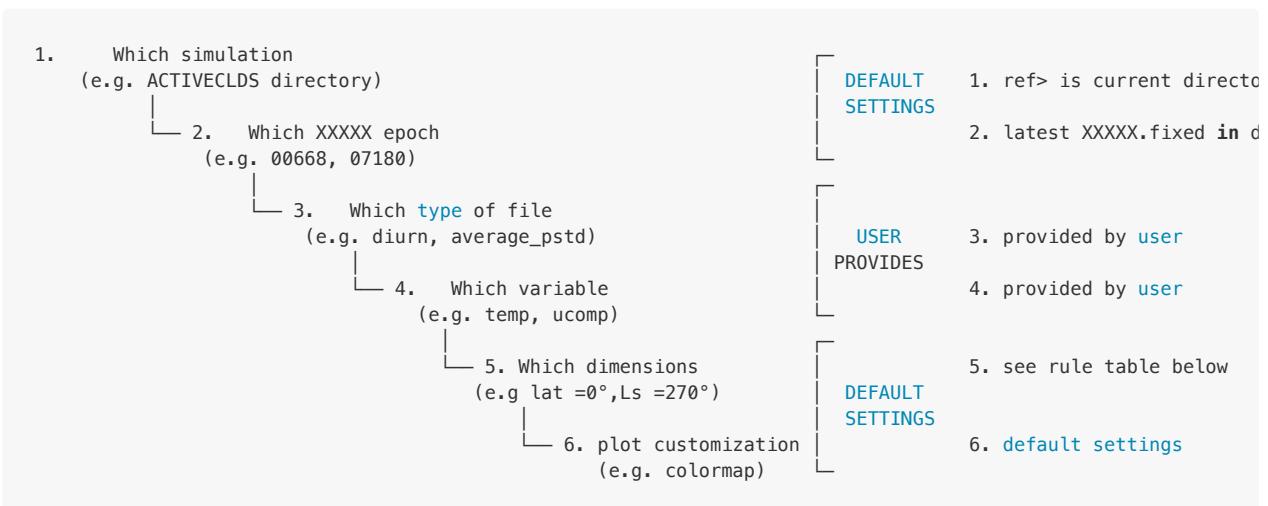
Let's remind ourselves that in order to create such plots from a **multi-dimensional** dataset, we first need to specify the **free** dimensions, meaning the ones that are **not** plotted.

## 3D data set



A refresher on cross-section for multi-dimensional datasets

The data selection process to make any particular cross section is shown in the decision tree below. If an effort to make the process of generating multiple plots as **streamlined** as possible, MarsPlot selects a number of default settings for the user.



The free dimensions are set by default using day-to-day decisions from a climate modeler's perspective:

| Free dimension | Statement for default setting                                                 | Implementation                    |
|----------------|-------------------------------------------------------------------------------|-----------------------------------|
| time           | "I am interested in the most recent events"                                   | time = Nt (last timestep)         |
| level          | "I am more interested in the surface than any other vertical layer"           | level = sfc                       |
| latitude       | "If I have to pick a particular latitude, I would rather look at the equator" | lat=0 (equator)                   |
| longitude      | "I am more interested in a zonal average than any particular longitude"       | lon=all (average over all values) |

| Free dimension | Statement for default setting                                                                               | Implementation |
|----------------|-------------------------------------------------------------------------------------------------------------|----------------|
| time of day    | "3pm =15hr Ok, this one is arbitrary. However if I use a diurn file, I have a specific time of day in mind" | tod=15         |

Rule table for the default setting of the free dimensions

In practice, these cases cover 99% of the work typically done so whenever a setting is left to default (= `None` in MarsPlot's syntax) this is what is being used. This allows to considerably streamline the data selection process.

`Custom.in` can be modified using your preferred text editor (and renamed to your liking). This is an example of the code snippet in `Custom.in` used to generate a lon/lat cross-section. Note that the heading is set to `= True`, so that plot is activated for MarsPlot to process.

```
<<<<<<<| Plot 2D lon X lat = True |>>>>>>>>
Title      = None
Main Variable = atmos_average.temp
Cmin, Cmax = None
Ls 0-360   = None
Level [Pa/m] = None
2nd Variable = None
Contours Var 2 = None
Axis Options : lon = [None,None] | lat = [None,None] | cmap = jet | scale = lin | proj = cart
```

In the example above, we are plotting the air temperature field `temp` from the `atmos_average.nc` file as a lon/lat map. `temp` is a 4D field (*time, level, lat, lon*) but since we left the time (`Ls 0-360`) and altitude (`Level [Pa/m]`) unspecified (i.e. set to `None`) MarsPlot will show us the *last timestep* in the file and the layer immediately adjacent to the *surface*. Similarly, MarsPlot will generate a *default title* for the figure with the variable's name (`temperature`), unit (`[K]`), selected dimensions (`last timestep, at the surface`), and makes educated choices for the range of the colormap, axis limits etc ... All those options are customizable, if desired. Finally, note the option of adding a secondary variable as **solid contours**. For example, one may set `2nd Variable = fixed.zsurf` to plot the topography (`zsurf`) from the matching `XXXXX.fixed.nc` file.

To wrap-up (the use of `{}` to overwrite default settings is discussed later on), the following two working expressions are strictly equivalent for `Main Variable = (shaded contours)` or `2nd Variable = (solid contours)` fields:

```
variable           SIMPLIFY TO           variable
|                   |                   |
00668.atmos_average@1.temp{lev=1000;ls=270} >>> atmos_average.temp
|       |       |       |       |
epoch  file type simulation    free dimensions   file type
|       |       |       |       |
file type simulation directory
```

These are the four type of accepted entries for the free dimensions:

| Accepted input                  | Meaning                                                                                                  | Example |
|---------------------------------|----------------------------------------------------------------------------------------------------------|---------|
| <code> --  - -- </code>         |                                                                                                          |         |
| <code>  None</code>             | <code>  Use default settings from the rule table below  Ls 0-360 = None  </code>                         |         |
| <code>  value</code>            | <code>  Return index closest to requested value in figure's unit   Level [Pa/m] = 50 (50 Pa) </code>     |         |
| <code>  Val Min, Val Max</code> | <code>  Return the averages between two values   Lon +/-180 = -30,30  </code>                            |         |
| <code>  all</code>              | <code>  all is a special keyword that return the average over all values in file   Latitude = all</code> |         |

Accepted values for the `Ls 0-360`, `Level [Pa/m]`, `Lon +/-180`, `Latitude` and `time of day` free dimensions

The time of day (`tod`) in diurn files is always specified using brackets `{}`, e.g.: `Main Variable = atmos_diurn.temp{tod=15,18}` for the average between 3pm and 6pm. This has allowed to streamlined all templates by

not including the *time of day* free dimension, which is specific to diurn files.

## MarsPlot.py: How to?

This section discusses MarsPlot capabilities. Note that a compact version of these instructions is present as comment at the very top of a new `Custom.in` and can be used as a quick reference:

```
===== |MarsPlot V3.2| =====
# QUICK REFERENCE:
# > Find the matching template for the desired plot type. Do not edit any labels left of any '=' sign
# > Duplicate/remove any of the <<< blocks>>>, skip by setting <<< block = False >>>
# > 'True', 'False' and 'None' are capitalized. Do not use quotes '' anywhere in this file
etc...
```

## Disable or add a new plot

Code blocks is set to `= True` instruct `MarsPlot` to draw those plots. Other templates in `Custom.in` are set to `= False` by default, which instructs `MarsPlot` to skip those plots. In total, `MarsPlot` is equipped to create seven plot types:

```
<<<<| Plot 2D lon X lat  = True |>>>>
<<<<| Plot 2D lon X time = True |>>>>
<<<<| Plot 2D lon X lev  = True |>>>>
<<<<| Plot 2D lat X lev = True |>>>>
<<<<| Plot 2D time X lat = True |>>>>
<<<<| Plot 2D time X lev = True |>>>>
<<<<| Plot 1D           = True |>>>> # Any 1D Plot Type (Dimension x Variable)
```

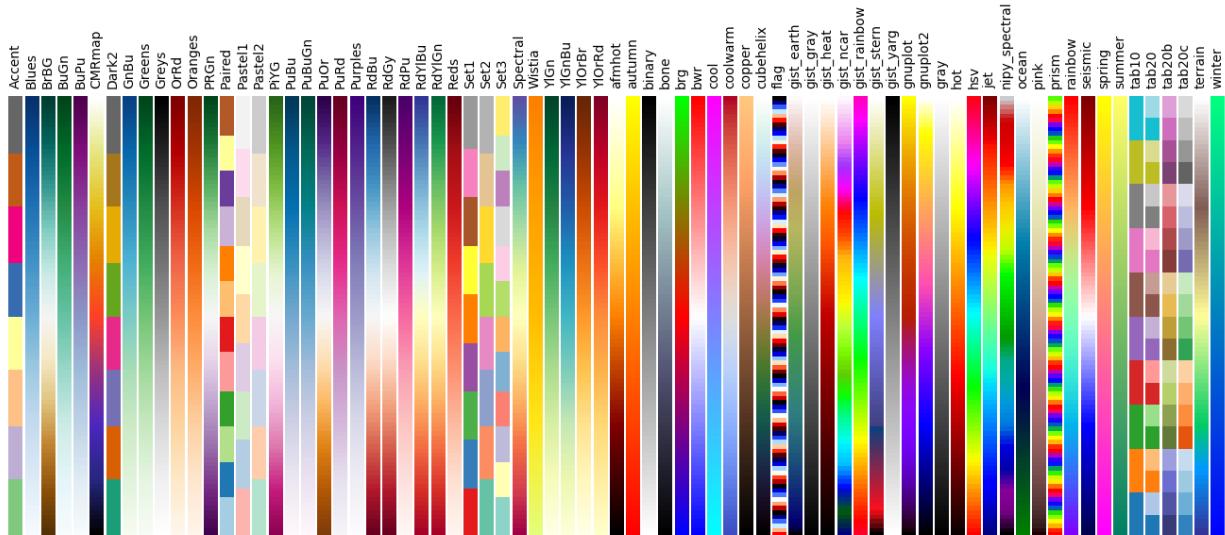
## Adjust the color range and colormap

`Cmin`, `Cmax` (and `Contours Var 2`) are how the contours are set for the shaded (and solid) contours. If only two values are included, MarsPlot use 24 contours spaced between the max and min values. If more than two values are provided, MarsPlot will use those individual contours.

```
Main Variable  = atmos_average.temp      # filename.variable *REQUIRED
Cmin, Cmax    = 240,290                   # Colorbar limits (minimum, maximum)
2nd Variable   = atmos_average.ucomp      # Overplot U winds
Contours Var 2 = -200,-100,100,200       # List of contours for 2nd Variable or CMIN, CMAX
Axis Options   : ls = [None,None] | lat = [None,None] | cmap = jet | scale = lin
```

Note the option of setting the contour spacing linearly `scale = lin` or logarithmically (`scale = log`) if the range of values spans multiple order of magnitudes.

The default colormap `cmap = jet` may be changed using any Matplotlib colormaps. A selection of those are listed below:



Finally, note the use of the `_r` suffix (reverse) to reverse the order of the colormaps listed in the figure above. From example, using `cmap = jet` would have colors spanning from `blue > red` and `cmap = jet_r red > blue` instead

*Supported colormap in Marsplot. The figure was generated using code from [the scipy webpage](#).*

## Make a 1D-plot

The 1D plot template is different from the others in a few key ways:

- Instead of `Title`, the template requires a `Legend`. When overplotting several 1D variables on top of one another, the `legend` option will label them instead of changing the plot title.
- There is an additional `linestyle` axis option for the 1D plot.
- There is also a `Diurnal` option. The `Diurnal` input can only be `None` or `AXIS`, since there is syntax for selecting a specific time of day using parenthesis (e.g. `atmos_diurn.temp{tod=15}`) The `AXIS` label tells `MarsPlot` which dimension serves as the X axis. `Main Variable` will dictate the Y axis.

Some plots like vertical profiles and latitude plots use instead Y as the primary axis and plot the variable on the X axis

```
<<<<<<<<<| Plot 1D = True |>>>>>>>>
Legend      = None          # Legend instead of Title
Main Variable = atmos_average.temp
Ls 0-360    = AXIS          # Any of these can be selected
Latitude     = None          # as the X axis dimension, and
Lon +/-180   = None          # the free dimensions can accept
Level [Pa/m] = None          # values as before. However,
Diurnal [hr] = None          # ** Diurnal can ONLY be AXIS or None **
```

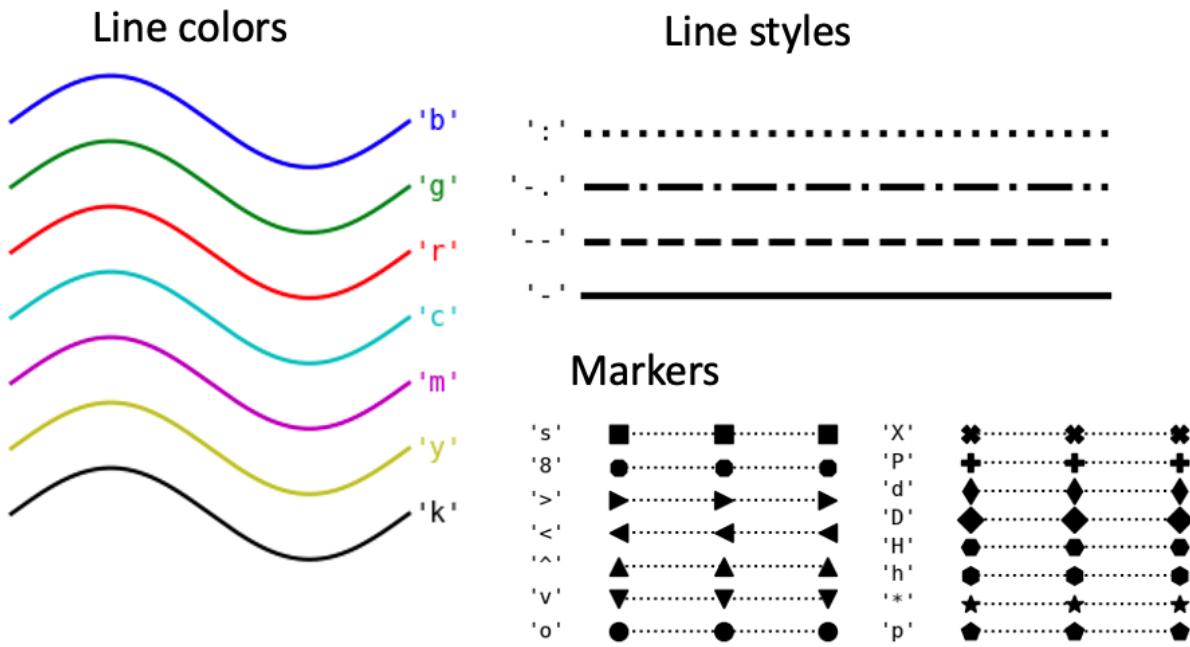
## Customize 1D plots

`Axis Options` specify the axes limits, and linestyle 1D-plot:

| 1D plot option                                       | Usage                                                  | Example                                             |
|------------------------------------------------------|--------------------------------------------------------|-----------------------------------------------------|
| <code>lat,lon+/-180,[Pa/m],sols = [None,None]</code> | range for X or Y axes limit depending on the plot type | <code>lat,lon+/-180,[Pa/m],sols = [1000,0.1]</code> |

| 1D plot option                  | Usage                                            | Example                                                           |
|---------------------------------|--------------------------------------------------|-------------------------------------------------------------------|
| <code>var = [None, None]</code> | range for the plotted variable on the other axis | <code>var = [120, 250]</code>                                     |
| <code>linestyle = -</code>      | Line style following matplotlib's convention     | <code>linestyle = -ob</code> (solid line & blue circular markers) |
| <code>axlabel = None</code>     | Change the default name for the axis             | <code>axlabel = New Temperature [K]</code>                        |

Here is a sample of colors, linestyles and marker styles that can be used in 1D-plots



*Supported colormap in Marsplot. This figure was also generated using code from [scipy-lectures.org](http://scipy-lectures.org)]*

## Put multiple plots on the same page

You can sandwich any number of plots between the `HOLD ON` and `HOLD OFF` keywords to group figures on the same page.

```
> HOLD ON
>
> <<<<| Plot 2D lon X lat = True |>>>>
> Title    = Surface C02 Ice (g/m2)
> ... (etc) ...
>
> <<<<| Plot 2D lon X lat = True |>>>>
> Title    = Surface Wind Speed (m/s)
> ... (etc) ...
>
> HOLD OFF
```

By default, MarsPlot will use a default layout for the plots, this can be modified by adding the desired number of lines and number of columns, separated by a comma: `HOLD ON 4 ,3` will organize the figure with a 4 -lines and 3-column layout.

Note that Custom.in comes with two plots pre-loaded on the same page.

## Put multiple 1D-plots on the same page

Similarly adding the `ADD LINE` keywords between two (or more) templates can be used to place multiple 1D plot on the same figure.

```
> <<<<<| Plot 1D = True |>>>>
> Main Variable      = var1
> ... (etc) ...
>
> ADD LINE
>
> <<<<<| Plot 1D = True |>>>>
> Main Variable      = var2
> ... (etc) ..
```

Note that if you combine `HOLD ON/HOLD OFF` and `ADD LINE` to create a 1D figure with several sub-plots on a **multi-figure page**, the 1D plot has to be the LAST (and only 1D-figure with sub-plots) on that page.

## Use a different epoch

If you have run a GCM simulation for a long time, you may have several files of the same type, e.g.:

```
00000.fixed.nc      00100.fixed.nc      00200.fixed.nc      00300.fixed.nc
00000.atmos_average.nc 00100.atmos_average.nc 00200.atmos_average.nc 00300.atmos_average.nc
```

By default MarsPlot counts the `fixed` files in the directory and run the analysis on the last set of files, `00300.fixed.nc` and `00300.atmos_average.nc` in our example. Even though you may specify the epoch for each plot (e.g. `Main Variable = 00200.atmos_average.temp` for the file starting at 200 sols), it is more convenient to leave the epoch out of the `Custom.in` and instead pass the `-date` argument to MarsPlot.

```
MarsPlot.py Custom.in -d 200
```

`-date` also accepts a range of sols, e.g. `MarsPlot.py Custom.in -d 100 300` which will run the plotting routine across multiple files.

There are several other plot customizations you can use:

- When two or more blocks are sandwiched between a `HOLD ON` and `HOLD OFF`, MarsPlot will draw the plots on the same page.
- Plots can be saved as images instead of PDFs by specifying your preferred filetype (PNG, EPS, etc.) when passing the `--output (-o)` argument to `MarsPlot`.
- When creating 1D plots of data spanning multiple years, you can overplot consecutive years by calling `--stack_year (-sy)` when submitting the template to `MarsPlot`.

## Access simulation in a different directory

The final plot-related functionality in `MarsPlot` is the simulation list, which allows you to point `MarsPlot` to different directories containing the MGCM output:

```
<<<<<<<<< Simulations >>>>>>>>>>>
ref> None
2> /path/to/another/sim                                # another simulation
3>
=====
```

Only 3 simulations have place holders but you can add additional ones if you would like (e.g. `4> ...` )

To access a variable from a file in another directory, just point to the correct simulation when calling `Main Variable` using the `@` character:

```
Main Variable = XXXXX.filename@N.variable`
```

Where `N` is the number in `<<< Simulations >>>` corresponding to the correct path.

## Overwrite the free dimensions.

By default, `MarsPlot` uses the free dimensions provided in each template (`Ls 0-360` and `Level [Pa/m]` in the example below) to reduce the data for both the `Main Variable` and the `2nd Variable`. You can overwrite this behavior by using parenthesis `{}`, containing a list of specific free dimensions separated by semi-colons ; The free dimensions within the `{}` parenthesis will ultimately be the last one selected. In the example below, `Main Variable` (shaded contours) will use a solar longitude of 270° and a pressure of 10 Pa, but the `2nd Variable` (solid contours) will use the average of solar longitudes between 90° and 180° and a pressure of 50 Pa.

```
<<<<<<<| Plot 2D lon X lat = True |>>>>>>>>
...
Main Variable = atmos_average
...
Ls 0-360      = 270
Level [Pa/m]   = 10
2nd Variable  = atmos_average{ls=90,180;lev=50}
```

## Element-wise operations

You can encompass variables between square brackets [] to perform element-wise operations, which is useful to compare simulations, apply scaling etc... `MarsPlot` will first load each variables encompassed with the brackets, and then apply the algebraic expression outside the [] before plotting.

These are examples of potential applications:

```
> Main Variable = [fixed.zsurf]/(10.**3)          (convert topography from [m] to [km])
> Main Variable = [atmos_average.taudust_IR]/[atmos_average.ps]*610 (normalize the dust opacity)
> Main Variable = [atmos_average.temp]-[atmos_average@2.temp]    (temp. difference between ref simu and
> Main Variable = [atmos_average.temp]-[atmos_average.temp{lev=10}] (temp. difference between the default (
```

## Commenting out and speed-up processing

Comments are preceded by `#`, following python's convention. Each `<<<|` block `|>>>` must stay integral so comments may be inserted between templates or comment all lines of the template (which is why it is generally easier to simply set the `<<<|` block = `False` `|>>>`) but not within a template.

You will notice the `START` key word at the very beginning of the template.

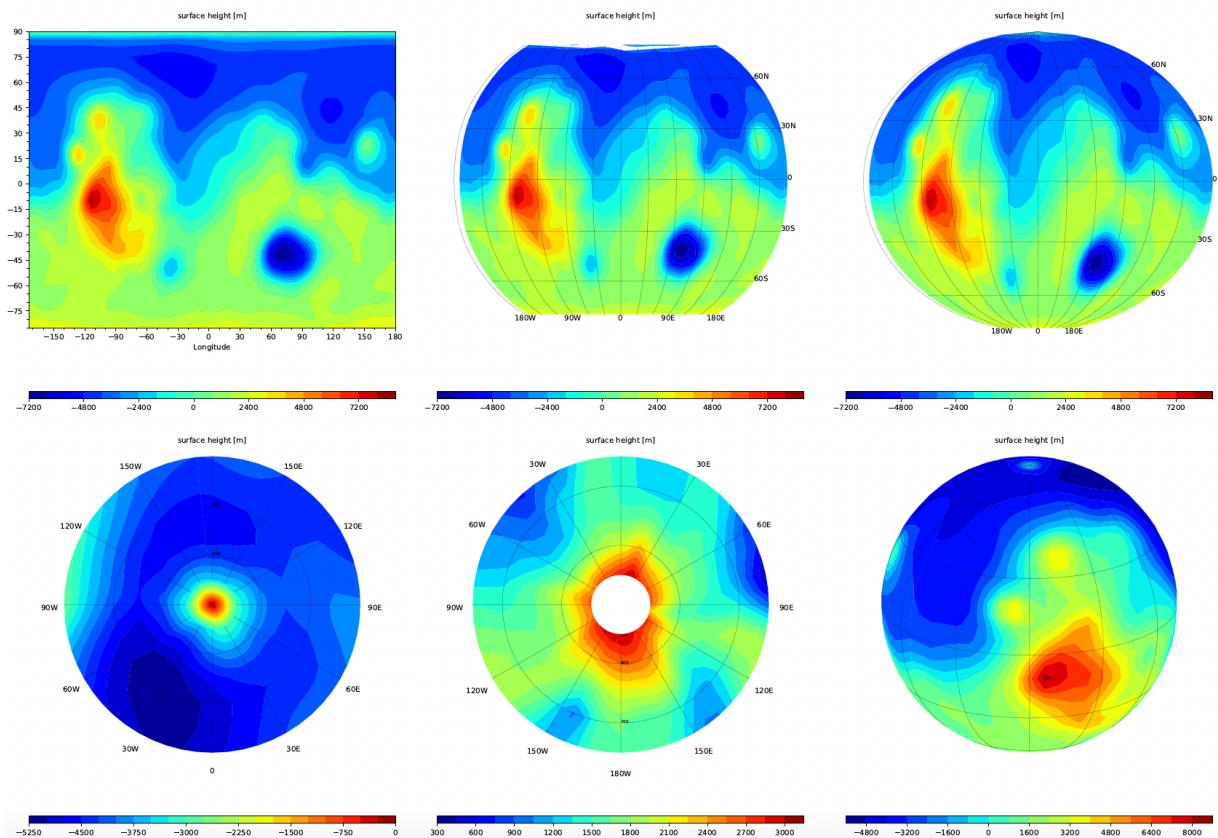
```
=====
START
```

This instructs MarsPlot to start parsing templates at this point. If you are already happy with multiple plots, you can move the `START` keyword further down in the Custom.in to skip those first plots instead of setting those to `<<<| Plot = False |>>>` individually. When you are done with your analysis, move `START` back to the top to generate a pdf with all the plots.

Similarly, you can use the keyword `ST0P` (which is not initially present in Custom.in) to stop the parsing of templates. In this case, the only plots processed would be the ones between `START` and `ST0P`.

## Change projections

For `Plot 2D lon X lat` figures, MarsPlot supports 3 types of cylindrical projections : `cart` (cartesian), `robin` (robinson), `moll` (mollweide), and 3 types of azimuthal projections: `Npole` (north polar), `Spole` (south polar) and `ortho` (orthographic).



(Top) cylindrical projection `cart` , `robin` and `moll` . (Bottom) azimuthal projections `Npole` , `Spole` and `ortho`

The azimuthal projections accept optional arguments as follows:

```
`Npole lat_max`  
`Spole lat_min`  
`ortho lon_center, lat_center`
```

## Figure format, size

As shown in the `--help` documentation of MarsPlot, the output format for the figure is chosen using the `--output (-o)` flag between `pdf` (default, requires the ghostscript software), `png`, or `eps`.

The `-pw` (pixel width) flag can be used to change the page width from its default value of 2000 pixels.

The `--vertical (-vert)` can be used to make the pages vertical instead of horizontal

## Access CAP libraries and make your own plots

CAP libraries are located (and documented) in `FV3_utils.py`. Spectral utilities are located in `Spectral_utils.py`, classes to parse fortran binaries and generate netCDF files are located in `Ncdf_wrapper.py`

The following code demonstrate how one can access CAP libraries and make plots for its own analysis:

```
#===== Import python packages =====
import numpy as np                                # for array operations
import matplotlib.pyplot as plt                     # python plotting library
from netCDF4 import Dataset                         # to read .nc files
#-----

# Open a fixed.nc file, read some variables and close it.
f_fixed=Dataset('/path_to_file/00000.fixed.nc','r')
lon=f_fixed.variables['lon'][:]
lat=f_fixed.variables['lat'][:]
zsurf=f_fixed.variables['zsurf'][:]
f_fixed.close()

# Open a dataset and read the 'variables' attribute from the NETCDF FILE
f_average_pstd=Dataset('/path_to_file/00000.atmos_average_pstd.nc','r')
vars_list      =f_average_pstd.variables.keys()
print('The variables in the atmos files are: ',vars_list)

# Read the 'shape' and 'units' attribute from the temperature VARIABLE
Nt,Nz,Ny,Nx = f_average_pstd.variables['temp'].shape
units_txt    = f_average_pstd.variables['temp'].units
print('The data dimensions are Nt,Nz,Ny,Nx=',Nt,Nz,Ny,Nx)
# Read the pressure, time, and the temperature for an equatorial cross section
pstd        = f_average_pstd.variables['pstd'][:]
areo        = f_average_pstd.variables['areo'][0] #solar longitude for the 1st timestep
temp        = f_average_pstd.variables['temp'][0,:,:18,:] #time, press, lat, lon
f_average_pstd.close()

# Get the latitude of the cross section.
lat_cross=lat[18]

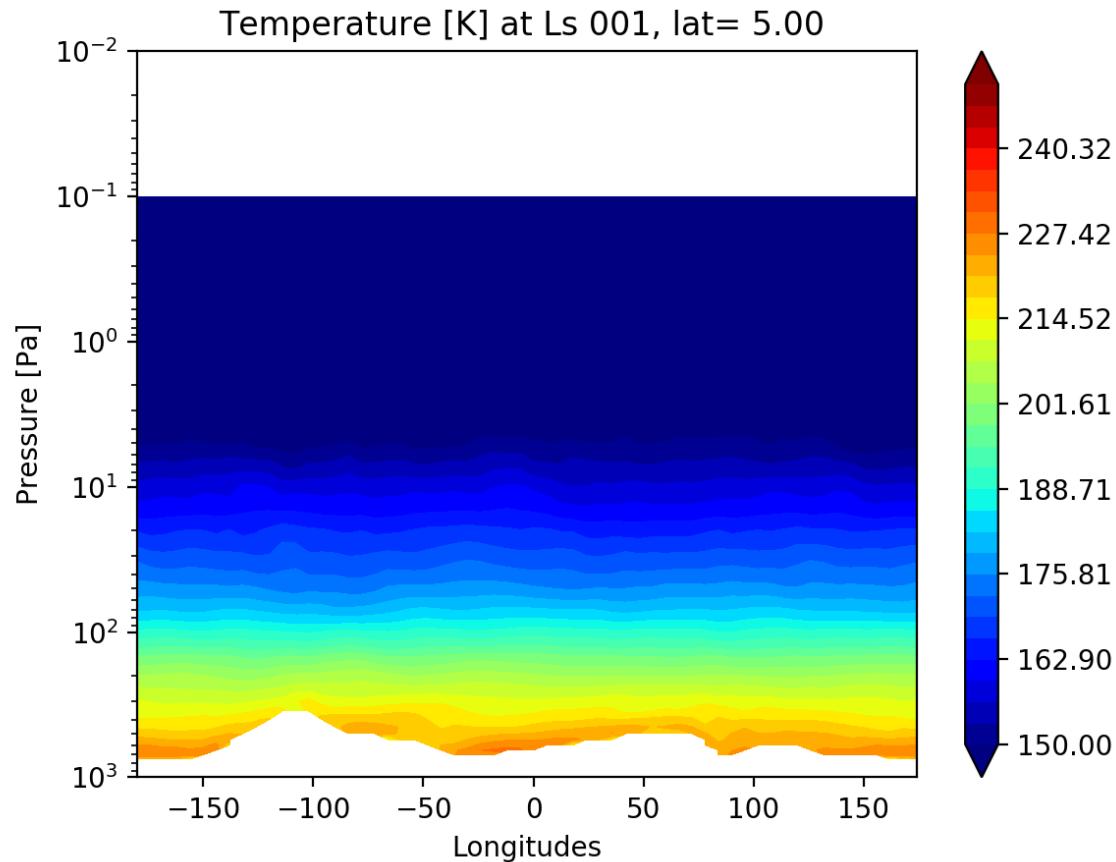
# Example of accessing functions from the Ames Pipeline if we wanted to plot
# the data in a different coordinate system (0>360 instead of +/-180 )
#-----
from amesgcm.FV3_utils import lon180_to_360,shiftgrid_180_to_360
lon360=lon180_to_360(lon)
temp360=shiftgrid_180_to_360(lon,temp)

# Define some contours for plotting
conts= np.linspace(150,250,32)

#Create a figure with the data
plt.close('all')
ax=plt.subplot(111)
plt.contourf(lon,pstd,temp,conts,cmap='jet',extend='both')
plt.colorbar()
```

```
# Axis labeling
ax.invert_yaxis()
ax.set_yscale("log")
plt.xlabel('Longitudes')
plt.ylabel('Pressure [Pa]')
plt.title('Temperature [%s] at Ls %03i, lat= %.2f'%(units_txt,areo,lat_cross))
plt.show()
```

will produce the following image:



## Debugging

`MarsPlot` is designed to make plotting MGCM output easier and faster so it handles missing data for you. For example, when dimensions are omitted with `None`, `MarsPlot` makes educated guesses for data selection and will tell you exactly how the data is being processed both in the title for the figures (if `Title = None`), and in the terminal output. Specifics about this behavior are detailed in the instructions at the top of `Custom.in`.

`MarsPlot` handles many errors by itself. It reports errors both in the terminal and in the generated figures. To by-pass this behavior (when debugging), use the `--debug` option with `MarsPlot` which will raise standard Python errors.

[Back to Top](#)