

# MESA tutorial: Session 2

## MESA output + Intermediate mass stars

### 1 Working with MESA output

The following is a summary of the official MESA documentation. For a more comprehensive overview of available output options, please refer directly to the documentation. You might have noticed that MESA stores its output as plain text files in the `LOGS` directory. There are two main types of output:

- **history.data:** logs each model's evolution in one line per model. The first line contains column indices, the second column names, and the remainder the data. Note: in case of a restart, older models are not removed—new data is appended. As a result *the model numbers are not guaranteed to be monotonically increasing!* Existing tools (like MESA Reader below) automatically take care of this, but if you write your own parser, be aware of this.
- Each **profile\*\*.data:** file contains a snapshot of a single model's structure. Each profile includes both global properties of the star, such as its age, and a large set of properties for each point in the model of the star given one line per point. In each case, the lines of data are preceded by a line with column numbers and a line with column names.

Lastly, you might also notice a **profiles.index** file. This maps profile filenames to model numbers. MESA saves profiles only at selected steps. Each line lists a model number, its priority (1 or 2), and the corresponding profile number. We will control some of the output of your MESA models in section 3.

**Using MESA Reader** MESA Reader is a Python framework with which one can easily plot data from the history and profile files. Read through the section on MESA Reader on the website to get familiar with its capabilities.

We want to install MESA Reader, but in such a way that we can always activate/access it, regardless of which computer we are using in the Faculty. Turns out that python **virtual environments** are exactly what we are looking for!

1. **Create a virtual environment** by running the following command in your terminal:

```
1 mkdir -p ~/venvs      # make dir to store venvs
2 cd ~/venvs            # go to that directory
3 python -m venv myenv   # create a venv called myenv
4 source ~/venvs/myenv/bin/activate # activate venv
```

You will notice that your terminal prompt changes, indicating that you are now working inside the virtual environment. To deactivate the virtual environment, simply run: `deactivate`.

2. Install Mesa Reader in your virtual environment by running:

```
1 pip install mesa_reader
```

You can install all kinds of useful python packages like this in your venv using `pip install <package_name>`. Though note that you can only install packages in your venv when it is activated!

Alternatively, you can directly git clone (download) the code from the project's Github repository somewhere in your **home** directory (e.g. `~/codes`)<sup>1</sup> by running:

```
1 git clone https://github.com/wmwolf/py_mesa_reader.git
```

Then run `python setup.py install` inside the cloned directory.

3. Throughout this course, we will use Jupyter Notebook to analyse our results. This is accessed via <https://jupyterhub.science.ru.nl/> and logging in with your science account. When you are logged in, activate your virtual environment by opening a terminal in Jupyter and running

```
1 source ~/venvs/myenv/bin/activate
```

Next, we need to register our venv as a kernel in the Jupyter hub, you only have to do this once. Run the following command in a *Jupyter Hub* terminal:

```
1 pip install ipykernel      # to run notebooks in your venv
2 python -m ipykernel install --user --name=myenv --display-name "
   Python (myenv)"
```

4. Open a new notebook in Jupyter by clicking on the plus sign in the top right, and from the new launcher select under Notebook: Python (myenv). In this notebook you should now be able to import MESA Reader without any errors:

```
1 import mesa_reader as mr
```

Note some other useful imports for calculating and plotting purposes could be:

```
1 import numpy as np
2 import matplotlib.pyplot as plt
3 import astropy.units as u
4 import astropy.constants as const
```

What `pip install` commands do you need to run to get all the aforementioned packages?

## 2 Overshooting in an intermediate-mass star

We will now investigate the effect of overshooting. In order to do this, we evolve an intermediate-mass star for different levels of convective overshooting.

1. **Download and set up the work folder.** Instead of starting from the beginning and copying the model work directory like last time to edit, a MESA model has largely been constructed for you already.
  - (a) Download the tar file from Brightspace and unpack it. This can be done using the archive manager by extracting the contents to your desired location on the **scratch** disc, or using the terminal by typing `tar -xvf session2.tar` in the directory where you downloaded the tar file and then moving the directory to your desired location on the **scratch** disc using a `cp -R` command. This file contains a MESA work folder in which we will work.

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<sup>1</sup>Note that the symbol `~` points to your home directory.

- (b) Once you open the `inlist_project` file, you will find a number of controls available for the model. First, choose a mass for the star between  $2.5 M_{\odot}$  and  $10 M_{\odot}$ , and modify the `inlist` accordingly. Note that the stopping condition allows the star to evolve up to the end of core helium burning. You may further notice several lines commented out concerning Convective overshooting. We will run several models for different values of the overshooting parameter.
- (c) First, compile (`./mk`) and run (`./rn`) the model without any overshooting. Use the various **PGSTAR** windows to follow and understand the evolution of this star. One of the **PGSTAR** windows is a Kippenhahn diagram, which shows information about the structure of the star as it evolves. Also try to understand this plot.
- (d) After the model finishes, create a new copy of the work folder and rename the **old** folder to an appropriate name (e.g. identify it by the mass and overshoot value used, like `M3_0ov0` for mass 3 and overshoot 0). In the new work folder, uncomment the lines relating to convective overshooting in your `inlist` and change the overshooting parameter `overshoot_f(1)` to 0.25. Look for the meaning of this and other overshooting parameters in the file `$MESA_DIR/star/defaults/controls.defaults` or by checking the documentation online.<sup>2</sup> Then compile and run the code again. Repeat this process for an overshooting parameter of 0.5.

#### Pro Tip

It is a good idea to organise your **MESA** work folders in a logical directory structure, in order to not make your home directory a mess. Just make sure you change the path in your **Python** script, such that it points to the right file.

2. **Inspecting the results.** We can now analyse the data in the 3 history files. For this import **MESA Reader** into your python notebook, as explained above.

To read in the data, you need to move the `history.data` files from your `scratch` directory to somewhere in your plotting folder, for example inside a folder called `data` and there inside a folder called after its run (e.g. `M3.0ov0`). For this example, you use the following line in your **Python** script:

```
1 f0_hist_data = mr.MesaData('data/M3.0ov0/LOGS/history.data')
```

To invoke the columns you want from the read-in data, in this case `f0_hist_data`, you use `f0_hist_data.X`, where `X` is the name of one of the columns of data inside `history.data`. The names of the columns can be found inside `history.data`, or you can print the available column names by typing:

```
1 print(f0_hist_data.bulk_names)
```

- (a) Make an HR diagram containing the 3 models. What changes do you see in the main-sequence evolution? What changes appear in the evolution *after* the main sequence? Can you explain these changes?
- (b) Make a plot of  $\rho_c$  vs  $T_c$ . How do the evolution tracks in this diagram change for different levels of overshooting?

<sup>2</sup>Note that the documentation online is for the latest version of **MESA**, which may differ slightly from the version you are using. The bottom right of the documentation page shows which version of the docs you are viewing, but note this only dates back to version r15140, which is when **MESA** was migrated to GitHub. `$MESA_DIR/star/defaults/controls.defaults` will always show the information for the correct version for your installation.

- (c) Construct a plot of the central helium abundance vs age for all 3 models and explain your findings.
- (d) By what fraction is the main sequence lifetime increased for an overshooting parameter of 0.25 compared to the model without overshooting? By what fraction does the *helium burning* lifetime change? Compare your findings with your neighbours who, hopefully, have chosen a star of different mass.

#### Pro Tip

As explained in section 2.2 of the first tutorial, it is a very good idea to copy your work folder from the `/scratch` directory to your home directory, after the MESA run has finished. This allows you to analyse your results using MESA Reader from any computer in the Faculty, not just the computer you ran your MESA models on!

## 3 Make your own Kippenhahn diagram

Kippenhahn diagrams (KHDs) show a star's internal structure over time and are useful for your final MESA project reports. Rather than taking screenshots from PGSTAR, you can use the Python script `mkipp`<sup>3</sup> to generate clearer, customizable plots. Here we will enable the necessary MESA output columns, help you set up `mkipp`, and get started with the scripts.

### 3.1 Setting MESA output

You can customize the MESA output through the `history_columns.list` and `profile_columns.list` files. In order to customize the output, just copy these files to your work directory, and uncomment the variables you want to include in the output. You can find default variations of these files in the `$MESA_DIR/star/defaults/` directory.

We will use the `mkipp` tool from <https://github.com/orlox/mkipp>. As you can read on the github page, to use `mkipp` you need to have the following MESA output available: “IMPORTANT!! your `history_columns.list` needs to have `mixing_regions` and `mix_relr_regions` specified for MESA to output the necessary data of the mixing regions in terms of mass and radius respectively. Also, newer versions of MESA include significantly less output in the profile files, so values such as `eps_nuc` need to be added to `profile_columns.list` as well.”

1. Pick one star that you like, and make sure that all these variables are uncommented. (i.e., remove the `!` in front of the variable) for both `history_columns.list` and `profile_columns.list` files of that star's work folder.
2. When you uncomment `mix_relr_regions`, it will most likely say `<integer>` after it. You should change this to 10, so that it has the same value as for `mixing_regions`.

For MESA to use the columns you have specified in these 2 files instead of the default set of columns, we have to make sure MESA knows that it needs to read these 2 files. To do this, we have to open `inlist_project` and include the following under `&star_job`:

```
1 ! to specify which output columns we want in history.data and
   profile.data
2 history_columns_file = 'history_columns.list'
3 profile_columns_file = 'profile_columns.list'
```

Now recompile and rerun your model.

<sup>3</sup><https://github.com/orlox/mkipp>

3. Next, actually cloning the `mkipp` repository:

```
1 cd ~/codes # or wherever you want to store the code
2 git clone git@github.com:orlox/mkipp.git
```

4. Unfortunately, `mkipp` is not available via `pip install`, so we have to make sure Python can find it. You can do this by adding the path to `mkipp` to your `PYTHONPATH` environment variable. To make your Jupyter notebook recognize and import a script or module from , you need to add that path to Python's search path at runtime.

```
1 import sys
2 import os
3 sys.path.append(os.path.expanduser("~/codes/mkipp"))
```

Your notebook will now look for python functions in the `/codes/mkipp` directory as well, and so you can import the three python modules that live in there:

```
1 import mkipp
2 import kipp_data
3 import mesa_data
```

5. To run a Kippenhahn diagram, you need to define the `logs_dir` argument of the `KippArgs` function as a list containing the path to your MESA output directory. For example, you can create a Kippenhahn diagram using the following command:

```
1 mkipp.kipp_plot(mkipp.Kipp_Args(logs_dirs = ['data/M3.0ov0/LOG']))
```

Do you understand what all the colors and hatching in the KHD mean?

6. Play around with the various options of `KippArgs` to customize your KHD. Have a look at the `mkipp/example.py` script for inspiration.