

WHAMP instructions/Advanced space plasma exercise

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1 Installing WHAMP

1. Get the WHAMP code by downloading at <https://github.com/ykempf/whamp> or using the git command

```
$ git clone git@github.com:ykempf/whamp.git whamp
```

which will download the code into the **whamp** directory.

2. Go to the code folder

```
$ cd whamp/src/
```

and compile the code

```
$ make
```

and do not worry about all the warnings. The file called **whamp** is the executable you will need.

By running this **whamp** executable you can use WHAMP on a point-by-point basis as the instructions in the manual tell you to do. But that is a bit painstaking so we use a script to automatise the process.

2 Running WHAMP from the Python scripts

1. Copy the **whamp** binary to the **exercise** directory.

```
$ cp whamp ../exercise
```

2. Run the script with Python

```
$ cd ../exercise
```

```
$ python3 <file>.py
```

and look at the 3D dispersion surfaces.

3. If you wish to save a png instead of displaying the result on screen, swap the **#** comment character between the last two lines of the script so that you use

```
matplotlib.pyplot.savefig('plot.png', dpi=DPI)
```

instead of

```
matplotlib.pyplot.show()
```

which will write the file **plot.png** to disk.