You may want to reprocess your spectroscopic data. This can be done on the **observer5** computer. The necessary software is installed in /opt/EffelsbergSpecPipeline. There is a bash script to set all the necessary paths. tcsh shell users should change to bash by typing bash. The teleskop, obs2 account are already bash accounts. Launch the following command:

# source /opt/EffelsbergSpecPipeline/init\_classwriter.sh

Now you should be ready to run the Offline Calibration pipeline. Type

### **OfflinePipeline**

Now you are in an ipython interface to reduce the data. There are the following commands to reduce the data:

## setFitsDir('/daten/Raw')

Set the source directory for the rawdata. The program looks in the directory for the raw mbfits files.

## setClassName('Filename','Directory')

this sets the output gildas, class file. if only:

# setClassName('Filename')

is given, the named file is created in the current dictionary. Without this command a file name e.g. like class\_2010\_10\_20.100m is created in the current directory. To write the files one has to know the scan number and must give the subscan number. Subscan numbers are 1 based. If the subscan is not in the file an error is plotted.

#### reduceSubscan(scannumber,subscannumber)

If you reduce the same scan twice, it is written twice to the class file with the same scan number. You can investigate the result with using\_class\_on\_observer3\_computer\_in\_effelsberg

## fixFFTSRefChannel(shiftFFTSRefChannel=True)

Before February 2011 the reference channel of the FFTS has to be shifted by +0.5 channels if the spectrum was not if flipped, otherwise -0.5 channels.

#### flipArrayOrientation(reverseArrayOrientation=True)

This function flips the dataArrays for the following reduce operations.

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