



BoA User Manual : APEX-MPI-MAN-0018

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# BoA – The Bolometer Data Analysis Software

## User and Reference Manual

### Purpose

The purpose of this document is to provide a description of the design and usage of the Bolometer Analysis (**BoA**) software package that was designed for the *Large APEX Bolometer Camera* (LABOCA) at **APEX**.

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v1.9	30.07.06	CV	all	version prepared for official approval
v1.10	01.08.06	CV	all	minor changes only
v1.11	02.08.06	CV	initial pages	Document history, Related documents, and Definintions added
v2.0	08.12.06	CV	all	new version created with different chapter structure plus updates. Significant changes to Chapters 3 (Cookbook) and 4 (Manual).
v2.1	15.12.06	CV	3	recipes for map, point and focus now present.
v2.2	01.03.07	CV	ch4	ch2 (install) also updated by RS
v2.3	08.03.07	CV	ch3 & 4	re-arrangement and updating
v2.4	09.03.07	CV	all	aesthetic changes to all plus updating of Ch4
v2.5	20.03.07	CV	ch4	brought up-to-date and re-organised

## Related documents

- RD-01** BoA User's manual
- RD-02** LABOCA design description, APEX-MPI-DSD-0016
- RD-03** Muders, Hafok, Wyrowski et al., 2006, A&A in press
- RD-04** The BoA Project: definition, F. Bertoldi et al. (June 2002)
- RD-05** A future bolometer data analysis software: requirements and definition, F. Bertoldi et al. (June 2002)
- RD-06** Initial BoA web site: <http://www.openboa.de>
- RD-07** LABOCA wiki: <http://www.astro.uni-bonn.de/abeelen/labocawiki>



## Definitions

For the following acronyms the understanding shall be:

<b>AIfA</b>	Argelander Institut für Astronomie der Universität Bonn
<b>AIRUB</b>	Astronomisches Institut der Ruhr-Universität Bochum
<b>APECS</b>	APEX Control Software
<b>APEX</b>	Atacama Pathfinder Experiment
<b>ASZCa</b>	APEX SZ Camera
<b>BoA</b>	Bolometer Array Analysis Package
<b>BoGLi</b>	BoA Graphics Library
<b>LABOCA</b>	Large APEX Bolometer Camera
<b>MAMBO</b>	Max-Planck Millimeter Bolometer
<b>MBfits</b>	Multi-beam fits format
<b>MPIfR</b>	Max-Planck-Institut für Radioastronomie, Bonn
<b>MOPSIC</b>	MAMBO data reduction software
<b>NIC</b>	IRAM bolometer reduction package
<b>SURF</b>	SCUBA data reduction software

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**Part I**

**User's Manual**

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# 1. INTRODUCTION

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The **Atacama Pathfinder Experiment (APEX)**<sup>1</sup> is a 12-meter radio telescope at the best accessible site for submillimeter observations, Llano de Chajnantor in Chile's Atacama desert.



Figure 1.0.1: The APEX telescope at Chajnantor in November 2003

LABOCA is a 295-channel facility bolometer camera for APEX. It operates in the 870  $\mu\text{m}$  atmospheric window and is to be commissioned in September 2006. It was built at the MPIfR bolometer lab by Dr. Ernst Kreysa and his staff.

**BoA** is a newly designed software package for the reading, handling, and analysis of bolometer array data. Its design and implementation is a collaborative effort of scientists at the MPIfR, AIfA and AIRUB that was started in 2002 and in part funded through a "Verbundforschung" grant to the MPIfR and RAIUB. **BoA** is an APEX facility software as part of the LABOCA instrument. The primary goal of **BoA** is to handle data from LABOCA at APEX, both for online visualization and offline processing. **BoA** could also be used to process data acquired with other instruments such as ASZCa at APEX or MAMBO at the IRAM 30-meter telescope. **BoA** includes most of the relevant functionalities of the current reduction packages (MOPSIC, NIC, SURF). The major difference is that **BoA** is written in a programming environment that is easier to modify, maintain, and re-use. Moreover, **BoA** naturally interfaces with APECS and the MBfits format.

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<sup>1</sup><http://www.mpifr-bonn.mpg.de/div/mm/apex/>

## 1.1 Philosophy and basic structure

### 1.1.1 Philosophy

BoA is designed with two major goals in mind: to provide a comprehensive tool for the reduction and analysis of data from the new generation of bolometer arrays, and to facilitate the extension and modification of the software by any user. **BoA** is intended to combine a simple and intuitive usage with the coverage of all aspects of data reduction, from raw data to final images. The natural choice for the creation of **BoA** is object oriented programming.

### 1.1.2 Programming language: Python

Most of **BoA** is written in Python, an interpreted, interactive and object-oriented programming language. Python does not adhere to all concepts of object-orientation as strictly as, e.g., C++ does. The resulting shortcomings can be overcome by sticking to some basic programming rules.

Python is a scripting language and as such allows **BoA** to be quickly and easily extended by the user. It also facilitates the wrapping of code written in C/C++ or FORTRAN. To improve execution speed, **BoA** computing-intensive tasks are therefore written in Fortran95.

### 1.1.3 Basic structure

**BoA** consists of a set of classes, most of which are defined in dedicated modules (files). In addition, a few functions are defined in separate modules. A detailed description of all classes and methods can be found in Sec. 3. The subdivision was chosen to reach a high modularity and an intuitive grouping of related functionalities within one class.

Two kinds of classes may be distinguished:

- **Data classes:** The `DataEntity` class defines the data structure which is used within **BoA**. Objects of this class contain the raw and reduced data and all relevant parameters of a single scan. This class also defines methods to fill the data object from an MBFITS file. Then, the `DataAna` class inherits from `DataEntity`: it contains all data related methods, plus some methods for data analysis (e.g. flagging, baseline). Then, the `Map` class inherits from `DataAna`: it contains all methods defined in `DataEntity` and `DataAna`, plus specific methods for map processing and display. Finally, classes dedicated to various observing modes inherit from the `Map` class: they contain additional methods specific to a given type of observation. Table 1.1 lists **BoA** data classes, with module names and short descriptions of their responsibilities.
- **Peripheral classes:** All other classes provide methods which either are used by data objects (e.g. `Image` is used within `Map` objects), or provide functionalities on the **BoA** level (e.g. `MessHand`). These classes are summarized in Table 1.2.

Finally, a few functions are defined in separate modules (listed in Table 1.3), which do not define any class. Thus, these functions can easily be imported and run from any level. In particular, the **BoAGraphic Library (BoGLi)** is defined in a collection of modules, which can be imported at the python level and do not require **BoA**. A description of **BoGLi** is given in Sect. 5.

Table 1.1: **BoA** data classes

class name	module	purpose
DataEntity	BoaDataEntity.py	data and parameters storage
DataAna	BoaDataAnalyser.py	general data analysis methods
Map	BoaMapping.py	map reduction
Focus	BoaFocus.py	focus reduction
Point	BoaPointing.py	pointing reduction
Sky	BoaSkydip.py	skydip reduction

Table 1.2: Other **BoA** classes

class name	module	purpose
Image	BoaMapping.py	image and axis description
Error	BoaError.py	
Help	BoaHelp.py	online help
MessHand	BoaMessageHandler.py	message handling
MamboMBFits	MamboMBFits.py	MAMBO to/from MB-Fits conversion
Timing	Utilities.py	benchmarking utilites

In addition, a number of utility and computing routines are written in Fortran modules. These routines are used within Python methods, and should in principle not be called directly by a **BoA** user.



Table 1.3: Other **BoA** modules

module name	purpose
<b>BoGLi</b> (see Sect. 5)	Graphic library
Utilities.py (see Sect. ??)	collection of utilities
BoaConfig.py (see Sect. ??)	global parameters definitions
BoaSimulation.py	LABOCA data simulator

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## 2. INSTALLATION

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This section describes how to install **BoA** and all required additional software packages, as well as how to update an existing **BoA** version.

### 2.1 Prerequisites

So far, **BoA** has been installed and tested on Mac OS X and on the following LINUX distributions:

- SuSE 10.0
- Scientific Linux 4.2
- Fedora Core 3
- Debian sid

The following software packages must be installed on a system to be able to install and run **BoA**. (The given version numbers indicate the versions that were used during development and tests with the respective LINUX distribution.)

#### 2.1.1 SuSE 10.0

Table 2.1: Prerequisites for SuSE 10.0

Package	Version
gcc / gcc-c++	4.0.2
compat-g77	3.3.5
readline-devel	5.0
libpng-devel	1.2.8
xorg-x11-devel	6.8.2
findutils-locate	4.2.23
subversion	1.2.3-2

Depending on the original setup, some or all of the packages specified in table 2.1 may already be installed on your system. Use SuSE's package manager *YaST* to check if the necessary packages are present and to install or update them if necessary.

### 2.1.2 Scientific Linux 4.2

Table 2.2: Prerequisites for Scientific Linux 4.2

Package	Version
gcc / gcc-c++	3.4.4
gcc-g77	3.4.4
readline-devel	4.3
libpng-devel	1.2.7
xorg-x11-devel	6.8.2
findutils	4.1.20
subversion	1.1.4-2

All packages listed in table 2.2 are part of a standard “Workstation” installation. If a package is missing or needs to be updated, use the package manager *yum* that comes with Scientific Linux.

### 2.1.3 Fedora Core 3

Table 2.3: Prerequisites for Fedora Core 3

Package	Version
gcc / gcc-c++	3.4.4
gcc-g77	3.4.4
readline-devel	4.3
libpng-devel	1.2.8
xorg-x11-devel	6.8.2
findutils	4.1.20
subversion	1.1.4-1.1

All packages listed in table 2.3 are part of a standard “Workstation” installation. If a package is missing or needs to be updated, use Fedora's package manager *yum*.

*NB:* Tests were performed with Fedora installed in a VMWare virtual machine. With its original kernel (version 2.6.9), Fedora run extremley slow. This is a know problem with Fedora Core 3 under VMWare. Updating to kernel version 2.6.12 improved the performance considerably.

Table 2.4: Prerequisites for Debian sid

Package	Version
gcc / gccc-c++	4.1.1-11
compat-g77	3.4.6-10
readline-devel	5.1-7
libpng-devel	1.2.8rel-5.2
xorg-x11-devel	1.0.0-8
findutils	4.2.28-1
subversion	

### 2.1.4 Debian sid

If a package listed in table 2.4 is missing or needs to be updated, use Debian's package manager *apt*. For any questions related with Debian, contact Alexandre Beelen.

### 2.1.5 Mac OS X

BoA has also been successfully installed and running on Apple Macintosh computers. Successful installations were done on a G4 laptop (Powerbook G4 550 MHz) and a G5 desktop (Powermac G5 2x1.8 GHz), under Mac OS X 10.3.5 to 10.3.9. BoA is also running on the new Intel Mac (installed and tested on a MacBook Pro 2.0 GHz, with Mac OS X 10.4.6). For any question related to **BoA** on Mac OS, please contact FredericSchuller (schuller@mpifr-bonn.mpg.de).

The prerequisites specified in table 2.5 refer to Mac OS X 10.4.6.

Table 2.5: Prerequisites for Mac OS X

Package	Version
XCode development tools	
python	2.3.5
gcc	4.0.1
gfortran	i386-apple-darwin8.7.1 gcc-4.2-20060805
fink toolkit	
swig	1.3.20
numeric-py23	23.1
readline	
libpng3	

## 2.2 Conflicts with other software

During the installation of **BoA** described in section 2.3, a set of programs and libraries (called *external packages*) that **BoA** needs is installed. This makes a complete installation of **BoA** more or less self-contained and reduces the chance of conflicts with other software installed on your system.

However, the behaviour of a system is not only determined by the set of software that is installed on it, but also by the environment that is defined both system-wide and on a per-user basis in various startup scripts. (For a complete list of the startup scripts consult the documentation of your system; most important are the startup scripts of the shell in use. See `man bash` and `man csh`)

During the installation of **BoA**, the installation script tries to set up an environment that allows a smooth installation. When running `boa`, **BoA**'s start-up scripts `.boarc.sh` and `.boarc.csh` try the same. However, there may be situations when this is not successful. If this is the case, careful inspection of the environment must be performed. (During the development phase of **BoA**, running **BoA** failed reproducibly on one particular system; after scrutinizing various startup scripts, the cause turned out to be a startup script for IRAF, that changed the C and Fortran compilers. After commenting out the IRAF related lines, **BoA** run without any further problems.)

## 2.3 Obtaining the installation script and packages

The command

```
svn checkout boa-install
```

will download the external packages and the installation script `install.sh` (written by Alexandre Beelen, Thomas Jürges, Frederic Schuller, and Reinhold Schaaf) from **BoA**'s Subversion repository to the directory `boa-install` in your current directory. The directory `boa-install` will be created if not already present.

You are now ready to start the installation. (The **BoA** software itself is not downloaded at this stage. It will be downloaded from the Subversion server during the installation.)

## 2.4 Installation using the `install.sh` script

Please note that the following instructions relate to the installation of **BoA** under Linux. For any question related to **BoA** on Mac OS, please contact FredericSchuller ([schuller@mpifr-bonn.mpg.de](mailto:schuller@mpifr-bonn.mpg.de)).

### 2.4.1 Running the `install.sh` script

Before running the `install.sh` installation script make sure that you have fulfilled the prerequisites described in Sect. 2.1!

1. Go to the directory where you have downloaded the `openboa` directory and files. Change into the directory `install` where the installation script `install.sh` is stored.

2. Run the *install.sh* script by typing:

```
./install.sh
```

This begins the process of installing **BoA**.

The script will prompt you for some paths (reasonable defaults are offered). If you don't want to use the default path, then please enter your chosen path, e.g. */home/smueller/BoA*, when prompted. (Please note that you have to specify the full path name explicitly: *~/BoA* will not work!)

The script will create in this installation directory some sub-directories, where all necessary files will be installed. The required disk space is about 200 MB.

You will also be prompted to enter yes (y) or no (n) for the installation of each software package. For a fresh installation, you should install every package included (even if, say Python, is already present on your system).

If you want to install all included packages, you can run the installation by

```
./install.sh -all
```

The install script will prompt you considerably less often.

Skipping installation of packages is useful if you resume an aborted installation (see below). If you wish to, you can try to see whether **BoA** works with your preinstalled versions of software; however, that is at your own risk!

3. After the installation is complete, type

```
source ~/.boarc.sh      (if you are working in bash)
or
source ~/.boarc.csh     (if you are working in csh).
```

4. You can now run **BoA** by typing *boa* at the prompt!

### 2.4.2 Details of the installation process

The installation consists of three stages, all of which are performed by the installation script:

- Installation of the external packages necessary for **BoA**
- Installation of BoA itself, including documentation and example FITS files
- Installation of BoA's initialization files *.boarc.sh* and *.boarc.csh*

Below you find a description of each of these three stages.

After the installation, you will find a installation log in *install/build.stat*. If the installation fails, the install script will tell you that something went wrong and give you a place where you can find information related with the failure. In addition, you can consult *install/build.stat* for information about the earlier steps of the installation.

**Installation of external packages**

The installation script will install the following external packages:

Table 2.6: External packages

Package	Version
Python	2.3.2
Numeric	23.1
numarray	0.9
swig	1.3.23
Intel FORTRAN	8.1
scipy_distutils	3.3_33.571
f2py	2.44.240_1892
pgplot	5.2
pPGPLOT	1.3
slalib	
pySLALIB	0.4
blas / lapack	
cfitsio	2.49
pCFITSIO	
BoA-FFTW-Numpy	1.0
mpfit	
wcslib	4.1
dchelper	
apexFitsWriter	
apexCalibrator	

The installation script prompts you for the location of the external packages. The default should allways be correct.

Next, you are prompted for the directory where **BoA** is to be installed. If this directory already exists, you must confirm that choice. (This case is necessary to resume an aborted installation or to update the **BoA** software itself. In all other cases, install to a new directory!)

The script then installs all external packages into this directory.

For some packages, (e.g. `scipy_obsutils`) you are prompted whether you want the package to be updated via CVS. This may not be necessary, so you can safely answer `n`. If you do update, the installation script provides you the necessary information (CVS login and password). Be aware that the CVS server may be slow or even down. If this is the case, you are prompted after a timeout of about 2 min whether you want to proceed without the CVS update. If you are nervous, cancel the installation with `Ctrl-C` and resume the installation (see below).

### Installation of BoA

When the installation of the external packages is complete, the **BoA** software itself is installed. Since it is not included in the `boa-install` download, it is downloaded from the BoA repository now. (Please be aware that you may have to use your own Subversion login and password here!)

The script prompts you for a directory, where **BoA** is to be installed. You can choose any accessible directory.

After the installation of the **BoA** software, the documentation and example FITS files are downloaded from the **BoA** repository and installed. Again you are prompted whether and where you want these features to be installed.

### Installation of BoA's initialization files

As last step, the script installs the initialization files `.boarc.sh` and `.boarc.csh` to your home directory. These scripts define a runtime environment for **BoA** (setting shell variables, paths, and aliases) for `bash` (`.boarc.sh`) and `csh` (`.boarc.csh`). Before running **BoA**, type

```
source ~/.boarc.sh          (if you are working in bash)
```

or

```
source ~/.boarc.csh        (if you are working in csh)
```

You may want to add this to your shell's startup script.

## 2.5 Resuming an incomplete installation

To resume an incomplete installation, run

```
install/install.sh
```

again. When prompted for the directory to which **BoA** is to be installed, specify the same directory as in the aborted installation. (Do this even if you will not install any external packages; the information is needed for the initialization files!)

You can then safely skip all installation steps, that were performed successfully in the last installation run.

Please be aware that you are prompted for the variable `PGPLOT_DIR` after skipping the installation of `pgplot`. A reasonable default is offered. However, if you want to use a pre-installed `pgplot`, you can specify this here.

## 2.6 Installation FAQ

### 2.6.1 BoA fails to start

- `ImportError: No module named fUtilities`



the fortran modules have not been compiled. Go to the fortran directory and type `make`

- `ImportError: libifport.so.x`

you dont have the fortran librarie in you `$LD_LIBRARY_PATH`, please source the `boarc.xx` file or check your installation.

### 2.6.2 I can't open a Graphical Device

- check the `pgplot` and `p_pgplot` installation
- if trying to output `png` files, make sure that `libpng` was present when compiling `pgplot`.

### 2.6.3 Reading a MBFits file fails

- check the `cfitsio` and `pcfitsio` installation
- check that the version of `MBFits.xml/$MBFITSEXML` you are using match the file you are trying to read

## 2.7 Updating BoA

Depending on the changes in **BoA** that make an update necessary (or desirable), an update of **BoA** only or a complete update (external packages and **BoA**) may be necessary. Unfortunately, presently there is no systematic way to find out whether a update of the external packages is necessary. The best choice may be first to try an update of **BoA** only, and if you experience problems when running **BoA** after that update, make a complete update of the external packages and **BoA**.

### Updating BoAonly

Examine the shell variable `BOA_HOME_BOA` that is set in the initialization files `~/.boarc.sh` and `~/.boarc.csh`, to find out, where **BoA** is installed. Move to this directory:

```
cd $BOA_HOME_BOA
```

Update **BoA** from the **BoA** repository by typing

```
svn update
```

Change to the `fortran` subdirectory and compile **BoA**'s Fortran modules:

```
cd fortran
make clean all
```

**Updating external packages and BoA**

If you experience problems when running **BoA** after updating **BoA** only, update the external packages as well. For that, follow the instructions in section [2.3](#) to obtain a new installation script and the external packages from **BoA**'s Subversion repository. Then follow section [2.5](#) to replace the external packages of your current installation that need to be updated. Do not forget to update **BoA** itself as described in the preceding section.

If this does not result in a working installation, do a fresh installation according to section [2.4](#), possibly into a new directory.

---

## 3. BoACOOKBOOK

---

This section describes how to start up **BoA** for the first time and describes some typical **BoA** sessions, including making a map and solving a pointing and focus. These example sessions are intended to act as recipes to allow the beginner or occasional user to get on air quickly. Users already familiar with the content of this cookbook can find more detailed information in Chapter 4.

### 3.1 Starting up BoA

The most common way to invoke **BoA** is to simply type `boa` at the command prompt. **BoA** then prints a welcome message providing version information and changes the prompt to the `boa>` prompt. (Note that you are nevertheless still in the interactive Python layer).

When **BoA** is initiated it imports a set of modules, instantiates the most essential objects and makes the respective methods available using the start script *BoaStart.py*.

In certain circumstances, more advanced users may wish to invoke **BoA** via a Python session. This can be done in the following ways:

- start a Python session and at the Python prompt type

```
>>> from BoaStart import *
```

- call Python in interactive mode (*-i*) with the file *BoaStart.py*

```
python -i /home/user/boa/BoaStart.py
```

where */home/user/* is the user path to the *boa* directory.

### 3.2 Getting started

```
open()                % 1
indir('/home/user/data/') % 2
ils()                 % 3
proj('projectID')     % 4
read('filename')      % 5
```

A typical **BoA** session will usually require a data file as input and a graphic output device. Command 1 opens the default graphics device (pgplot). Command 2 sets the desired input directory, i.e. in this

case the input data file is located in a directory called `/home/user/data/`. The content of this directory can then be listed (command 3). The project ID can also be set (command 4) so that filenames may subsequently be described by just the observation number. Command 5 then reads in the input data file.

Note, these commands and those used in the sections below are abbreviations for the full user method names, as is described in Section 4.2.1.

### 3.3 Ending a session

To end a session you can first close the graphics device by typing

```
close()
```

then end the **BoA** session by typing `ctrl+d`.

### 3.4 Getting Help

You can get help on a **BoA** `command()` at any time by typing

```
print command.__doc__
```

at the prompt.

### 3.5 Example of making a map

```
open() % 1
indir('/home/user/data/') % 2
ils() % 3
proj('T-77.F-0002-2006') % 4
read('59491') % 5
signal() % 6
signal(1) % 7
mapping() % 8
medianBaseline() % 9
plotRmsChan() % 10
flagRms(above=1) % 11
flagRms(below=0.2) % 12
updateRCP('jup-44830-32-improved.rcp') % 13
flagPos(radius=150.) % 14
base(order=1) % 15
medianNoiseRemoval() % 16
plotRmsChan() % 17
flagRms(above=0.5) % 18
plotRmsChan() % 19
flagC([140,227]) % 20
```

---

```

despike() % 21
computeWeight() % 22
unflag(flag=5) % 23
mapping() % 24
doMap(system='EQ',sizeX=[83.9,83.73],sizeY=[-5.48,-5.28],oversamp=5.) % 25
smooth(6./3600.) % 26
mapdisp(caption=data.ScanParam.caption()) % 27
close() % 28

```

### 3.5.1 Setting up and accessing the data

The initial steps for starting up a typical session were described in Section 3.2. Command 1 opens the default graphics device and Command 2 sets the desired input directory. The content of this directory is then listed (command 3). The project ID can then be set (command 4) so that filenames may subsequently be described by just the observation number (in this example the file-naming convention is for LABOCA data, and consists of the APEX project ID (*T-77.F-0002-2006*) and the observation number). Command 5 then reads in the input data file for observation *59491*.

### 3.5.2 Visualising the data

To get a first look at the data you can use command 6 to plot the time series of the data for each pixel, or command 7 to look at the time series of the data for an individual pixel. You can also make a rough map using command 8. These commands will be used again (see below) when the data is processed.

### 3.5.3 Basic Processing and Analysis

Usually the processing of the data will begin with subtracting a baseline. This is done with command 9, where the median value per channel and per subscan is removed. To see the data after baseline subtraction you can use commands 6 and 7 again. Next, command 10 plots the RMS value versus pixel (channel) number. Commands 11 and 12 then flag pixels with RMS values which are higher or lower than the given value as bad. At this point you can use command 10 again to view the remaining unflagged data.

Command 13 reads in the rcp file for calibration. Command 14 then flags the area in which source signal is present, and commands 15 and 16 remove a baseline (using a polynomial fit) and the median noise value. The new distribution is then checked with command 19.

Bad channels can be flagged using command 20, and the data then despiked (command 21). If necessary, the despiked data can be examined using commands 6 and 7. Before producing a map the data should be weighted, in this case each channel weighted as the inverse of the square root of the rms of that channel (command 22). Command 23 then unflags the previously flagged source position.

There are then a number of ways to produce a final map. Command 24 produces a quick map in Az-El coordinates. Alternatively command 25 produces a map in EQ coordinates. See Chapter 4 for optional arguments for these and other methods. The map may then be smoothed (command 26) and this smoothed map displayed (command 27).

### 3.6 Example of solving a pointing

```

open() % 1
indir('/home/user/data/') % 2
proj('T-77.F-0002-2006') % 3
read('42947') % 4
signal() % 5
signal(1) % 6
mapping() % 7
medianBaseline() % 8
mapping(oversamp=3) % 9
solvepointing(plot=1) % 10
clear() % 11
read('46117') % 12
medianBaseline() % 13
medianNoiseRemoval() % 14
plotRmsChan() % 15
flagRms(above=20) % 16
mapping(oversamp=3) % 17
solvepointing(plot=1) % 18
close() % 19

```

The above recipe shows a typical session to solve a pointing. As usual (see Section 3.2) we begin by opening a graphics display device, setting the input directory, and setting the project ID (commands 1,2 and 3). The data file containing the pointing observation is read in (command 4), in this case a strong pointing source (Jupiter). As a first look at the data, the time series of the data for each pixel (command 5) or individual pixel (command 6) can then be plotted. Likewise a rough first-look map can be made (command 7). To construct the map on which to solve the pointing, the median baseline is first removed (command 8) (to see how the signal looks now you can repeat commands 5 and 6). Finally the pointing map is constructed (command 9) and the pointing solved (command 10).

If the pointing source is fainter (in this case an observation of Uranus), some additional steps could be taken. Following the above example, a graphics display window is already open so we can clear the display using command 11. Command 12 then reads in the data file containing the observation of Uranus. Again, first looks at the data can be made using commands 5, 6 and 7. The median baseline is then removed (command 13), and this time the median noise value is also removed (command 14). You can then check at what RMS most channels are using command 15, then use command 16 to flag channels with RMS values above a certain value (in this case an RMS of 20). Command 15 can then be repeated to see how the data looks now. The pointing map can then be constructed (command 18) and the pointing solved (command 18). Command 19 then closes the graphics display device.

### 3.7 Example of solving a focus

```

open() % 1
indir('/home/user/data/') % 2
proj('T-77.F-0002-2006') % 3
read('43275') % 4
solveFocus() % 5

```

```
read('46118')           % 6
medianBaseline()        % 7
medianNoiseRemoval()    % 8
solveFocus()            % 9
close()                 % 10
```

The above recipe shows a typical session to solve a focus. As usual (see Section 3.2) we begin by opening a graphics display device, setting the input directory, and setting the project ID (commands 1,2 and 3). Command 4 reads in the data file, in this case for a strong source (Jupiter). Command 5 then solves the focus. Command 6 then reads in a new data file, this time for a fainter source (Uranus). This time the median baseline and median noise levels are removed before solving the focus (commands 7, 8 and 9).

### 3.8 Demonstration scripts

The recipes described in Sections 3.5, 3.6 and 3.7 form the basis of three example scripts which are provided in order to demonstrate some of the basic functionalities of **BoA**. These are: *ExampleMap.py*, *ExamplePointing.py* and *ExampleFocus.py* and can be found in the directory */home/user/boa/examples/*. Run the scripts by typing:

```
execfile('/home/user/boa/examples/ExampleMap.py')
```

---

## 4. BoAUSER MANUAL

---

In this chapter you will find information about the structure of **BoA**, how **BoA** can be used, together with detailed descriptions of user methods. Since many user methods have an abbreviated form, these are listed in Section [4.18](#).

### 4.1 About BoA

In this Section we give a basic overview of the structure of **BoA**. Section [4.1.1](#) gives a brief introduction to the raw data file format, and Section [4.1.2](#) shows an overview of the data structure within **BoA**. More in-depth descriptions are given in Chapter [6](#).

#### 4.1.1 Input data

The data acquired at the APEX telescope are stored in a new file format, known as the MB-Fits format (for Multi-Beam FITS format, see the reference document APEX-MPI-IFD-0002 by Hatchell et al. for details). These files contain:

- the raw data as provided by the Frontend-Backend in use at the telescope
- data associated parameters: time of the observations, positions on the sky...
- a description of the complete Scan (eg. for a map: number of lines, steps between lines...)
- parameters of the receiver channels in the array: relative positions, relative gains

A more complete description of the input data format is given in Sect. [6.1](#).

#### 4.1.2 Internal data handling

Taking full advantage of the object-oriented nature of Python, **BoA** handles data by means of objects of various classes. The primary class for data storage and manipulation is called `DataEntity` (see also Section [??](#)). This class allows to store the raw data and associated parameters, and it provides methods relevant for any kind of observations (e.g. reading data from an MB-FITS file, plotting the signal as time series, plotting the telescope pattern). The most important attributes of this class are:

- `BolometerArray`: here, the relative positions and gains of the receiver channels are stored, as well as generic informations about the instrument and telescope (name, diameter, coordinates...)



- **ScanParam**: this contains the data associated parameters: coordinates of each point in several systems, timestamps (in LST and MJD), subscans related informations
- **Data**: this is a 2D array (time  $\times$  bolometer) which contains the current version of the data. At time of reading, the raw data are stored there; the content of this array is then altered by any processing step
- **DataFlags**, **DataWeights**: 2D arrays, with same size as **Data**, where flagging values and relative weights are stored for each individual data point

For processing different types of observations, **BoA** then provides several classes which inherits from **DataEntity**. Inheritance allows to define a class which contains all attributes and methods of the parent class, plus some specific attributes/methods. The inheritance scheme in **BoA** is as follows:

```
DataEntity < DataAna < Map < Point < Focus
```

When **BoA** is started, one object of class *Focus* is created with name *data*; this is the current data object, on which all reduction procedures can be applied. Additional objects of any data class can be created by the user within one **BoA** session. Then, applying processing methods to a data object with a different name than *data* requires to enter the full syntax (see Chapter ...), including the full name of the method, as opposed to the shortcuts described in Chapters 3 and 4.

**Note:** Python ensures no real difference between private and public attributes. There are only hidden attributes but this hiding can be overcome easily. Therefore the user might set any attribute directly and call any method. This is not advisable and may easily corrupt the whole **BoA** session. It is more recommendable to just use those methods for which the start script *BoaStart.py* provides abbreviations.

## 4.2 BoA usage

### 4.2.1 Methods

**BoA** tasks are accessed by directly calling the appropriate methods from the interactive Python layer. This ensures the full availability of all Python and ppgplot facilities. As the method names to be called from the Python layer may be rather long, the start script *BoaStart.py* provides a set of convenient abbreviations for those methods which are meant to be called directly by the user (“public” methods). We will therefore refer to these as user methods, a full list of which can be found in Section 4.18.

#### Example:

The name of the method to open a new graphic device is *DeviceHandler.openDev* and it can be called by

```
DeviceHandler.openDev()
```

or more conveniently by the abbreviations (user methods)

```
open() or op()
```

(note that the parentheses are always mandatory).

### 4.2.2 Arguments

Nearly all user methods require arguments to be passed. Nevertheless, the methods provide default arguments which thus may be omitted. In this case many methods just supply status information.

**Example:**

The user method `indir()` sets the desired input directory and requires the directory name as its argument:

```
indir('/home/user/data/')
```

The directory name is a string argument and has to be passed embedded in double or single quotes. Note that for consistency, in the examples throughout this manual we always use single quotes, but these can of course be substituted for double quotes.

Omitting the argument does not change the input directory but instead results in the supply of the current directory name:

```
indir()
```

In case an argument has to be typed more often a Python variable can be used:

```
a='/home/user/data/'  
indir(a)
```

Some methods require a list as argument. In Python a list is embedded in square brackets with a comma as separator. Python provides a variety of functionalities to manipulate lists.

**Example:**

The user method `signal()` plots the time series of the data (flux density or counts versus time). It allows the user to define the list of channels plotted:

```
signal([18,19,20])
```

To create a list you can use the Python function `range()`:

```
mylist=range(1,163)  
signal(mylist)
```

or:

```
signal(range(1,163))
```

Even if the list contains only one element the square brackets are mandatory:

```
signal([5])
```

User methods can also be called using keyword arguments of the form *keyword = value*.

**Example:**

By default, the user method `signal()` plots the signal versus time connecting the datapoints with lines:

```
signal()
```

However, if you prefer, for example, to see the individual datapoints without lines, you can modify the value of the *style* argument:

```
signal(style='p')
```

A description of graphics related arguments such as *style* is given in Section ??.

**4.2.3 Output**

Most user methods supply status information as screen output when being called. The amount of information displayed can be restricted using the message handler associated with the main *data* object:

```
data.MessHand.setMaxWeight(4)
```

where the argument is an integer value between 1 and 5, with the following meaning:

- 1: errors, queries
- 2: warnings
- 3: short info
- 4: extended info
- 5: debug

**4.3 Making maps**

Several methods are provided for constructing a map.

**4.3.1 Making a map**

METHOD: `doMap` (*optional arguments*)

DESCRIPTION: construct a map in EQ or (Az,El) coordinates

OPTIONAL ARGUMENTS:

---

<i>chanList</i>	channels to consider
<i>flag</i>	flag values to consider
<i>oversamp</i>	oversampling factor (beam fwhm / pixel size). Default=2.
<i>beammap</i>	compute a beam map (default: no)
<i>system</i>	coord. system, one of 'HO' (Az,El *offsets*) or 'EQ' (RA, Dec absolute coord.) default = 'HO'; optionally 'EQFAST' to do only one rotation on small maps (faster)
<i>sizeX</i>	limits in Az of the map
<i>sizeY</i>	limits in El of the map
<i>noNan</i>	remove NaN in self.Results?
<i>style</i>	color table to use in image
<i>smooth</i>	do we smooth with beam? (default: no)
<i>noPlot</i>	do not plot the map? (default: no)
<i>caption</i>	plot caption
<i>aspect</i>	keep aspect ratio? (default: yes)
<i>showRms</i>	compute and print rms/beam? (default: yes)
<i>rmsKappa</i>	kappa in kappa-sigma clipping used to compute rms

### 4.3.2 Making an Az-El map

METHOD: `horizontalMap` (*optional arguments*)

DESCRIPTION: construct a map in (Az,El) coordinates

OPTIONAL ARGUMENTS:

<i>chanList</i>	channels to consider
<i>flag</i>	flag values to consider
<i>oversamp</i>	oversampling factor (beam fwhm / pixel size). Default=2.
<i>sizeX</i>	limits in Az of the map
<i>sizeY</i>	limits in El of the map
<i>noNan</i>	remove NaN in self.Results?
<i>style</i>	color table to use in image
<i>noPlot</i>	do not plot the map? (default: no)
<i>caption</i>	plot caption

## 4.4 User methods for flagging data

### 4.4.1 Despiking

METHOD: `despike` (*optional arguments*)

DESCRIPTION: Flag yet unflagged data below 'below'\*rms and above 'above'\*rms.

OPTIONAL ARGUMENTS:

*chanList* list of channels to be flagged (default: current list)  
*below* flag data with value < 'below'\*rms  
*above* flag data with value > 'above'\*rms

#### 4.4.2 Flagging a list of channels

METHOD: `flagChannels` (*optional arguments*)

DESCRIPTION: assign flags to a list of channels. To unflag a channel simply flag with flag=0.

OPTIONAL ARGUMENTS:

*chanList* list of channels to be flagged (default: current list)  
*flag* flag value

#### 4.4.3 Flagging data by time interval

METHOD: `flagMJD` (*optional arguments*)

DESCRIPTION: flag data by MJD interval

OPTIONAL ARGUMENTS:

*channel* list of channels to flag (default: 'all')  
*below* flag data below this value (default end of the scan)  
*above* flag data above this value (default start of the scan)  
*flag* flag value to set (default 1)

#### 4.4.4 Flagging a position on the sky

METHOD: `flagPosition` (*optional arguments*)

DESCRIPTION: flag a position in the sky within a given radius

OPTIONAL ARGUMENTS:

*chanList* list of channels to flag (default: 'all')  
*Az/El* the horizontal reference position (arcsec for offsets, deg for absolute)  
*radius* aperture to flag in unit of the reference position  
*flag* flag to be set (default 5)  
*offset* flag on the offsets (default yes)

#### 4.4.5 Flagging channels with certain rms values

METHOD: `flagRms` (*optional arguments*)

DESCRIPTION: flag channels with rms below 'below' or above 'above'.

OPTIONAL ARGUMENTS:

*chanList* list of channel to flag (default: current list)  
*below* flag channels with rms < 'below'  
*above* flag channels with rms > 'above'  
*flag* flag value to set

#### 4.4.6 Flagging subscans

METHOD: `flagSubscan (optional arguments)`

DESCRIPTION: flag a list of subscans

OPTIONAL ARGUMENTS:

*subList* list of subscan numbers (or single number) to be flagged  
*flag* flag value to be set

#### 4.4.7 Unflagging data

METHOD: `unflag (optional arguments)`

DESCRIPTION: Unflag data, i.e. reset to 0.

OPTIONAL ARGUMENTS:

*channel* list of channels to be unflagged (default: current list)  
*flag* unflag only this value (default 1)

#### 4.4.8 Unflagging a list of channels

METHOD: `unflagChannels (optional arguments)`

DESCRIPTION: Unflag a list of channels

OPTIONAL ARGUMENTS:

*channel* list of channels to be unflagged (default: current list)

### 4.5 Baseline subtraction, sky removal and statistics

#### 4.5.1 Computing the Rms in a map

METHOD: `computeRms () (optional arguments)`

DESCRIPTION: compute rms/beam in a map

OPTIONAL ARGUMENTS:

*cell*            size of one beam in pixel  
*rmsKappa*    for kappa-sigma clipping before computing rms

#### 4.5.2 Computing weights

METHOD: `computeWeight()` (*optional argument*)

DESCRIPTION: compute weights and store them in `DataWeights` attribute

OPTIONAL ARGUMENTS:

*method*    type of weighting (default='rms', i.e. use  $1/\text{rms}^2$ )

#### 4.5.3 Median baseline removal

METHOD: `medianBaseline()` (*optional arguments*)

DESCRIPTION: remove median value per channel and per subscan

OPTIONAL ARGUMENTS:

*chanList*    list of channels to process (default: all; [] : current list)  
*subscan*    compute baseline per subscan (default: yes)

#### 4.5.4 Noise removal

METHOD: `medianNoiseRemoval()` (*optional arguments*)

DESCRIPTION: remove median noise from the data.

OPTIONAL ARGUMENTS:

*chanList*    list of channel to flag (default: all; [] : current list)  
*chanRef*    reference channel number (default: `RefChannel`)  
*computeFF*    compute skynoise FF (default) or use existing `FF_Median?`

#### 4.5.5 Polynomial baseline removal

METHOD: `polynomialBaseline()` (*optional arguments*)

DESCRIPTION: perform polynomial baseline removal on the data

OPTIONAL ARGUMENTS:

*chanList*    list of channels to flag (default: all; [] : current list)  
*order*    polynomial order, >0  
*subscan*    compute baseline per subscan (default: yes)  
*plot*    plot the signal and the fitted polynomials (default: no)  
*subtract*    subtract the polynomial from the data (default: yes)

### 4.5.6 Smoothing an image

METHOD: `smoothBy` (*optional arguments*)

DESCRIPTION: smooth the image with a 2D gaussian of given FWHM

OPTIONAL ARGUMENTS:

*beamSize* the FWHM of the smoothing gaussian

### 4.5.7 Obtaining the statistics

METHOD: `statistics`

DESCRIPTION: compute mean, median, rms for all scans and subscans for all used channels

## 4.6 Pointing

### 4.6.1 Solving a pointing

METHOD: `solvePointingOnMap` (*optional arguments*)

DESCRIPTION: compute the offset on the data.Map object

OPTIONAL ARGUMENTS:

*gradient* shall we fit a gradient ? (default: no)

*circular* fit a circular gaussian instead of an elliptical gaussian

*radius* use only bolo inside this radius (negative means multiple of beam) (default: 10 beams)

*Xpos* source position if using fixed position

*Ypos* source position if using fixed position

*fixedPos* if set, don't fit position, but use Xpos, Ypos

*plot* do we plot the results? (default: no)

*display* display the result of the fit (default: yes)

WARNING : No Smoothing should be applied to the map before using this function, or the fitted fwhm will be useless, use `fine oversamp` to make reasonable fit.

## 4.7 Focus

The recommended way to conduct Laboca focus observations is to perform a series of  $n \times 3$  short, symmetric on-offs, e.g. 3 or  $6 \times (4 \times 5 \text{sec})$ . For this simply the onoff has to be reduced and then the results can be fitted by a parabola.



### 4.7.1 Solving a focus

METHOD: `solveFocus`

DESCRIPTION: compute the optimal focus position

## 4.8 File reading

### 4.8.1 Reading a FITS file

Reading a FITS file into **BoA** is done with the `read()` command. You may want to define the input directory first:

```
indir('../fits/')      # set the input directory
read('filename')      # read file filename.fits
```

The data are then stored in the default *data* object. It is possible to use several data objects, and to store the content of a file to a user defined object requires the following syntax:

```
data2 = BoaMapping.Map() # define a second data object
                        # of class Map
data2.read('filename')
```

## 4.9 Controlling graphics display devices

In order to display your data in various ways using the **BoA** plotting methods described in Section 4.10 below, you first need to open a graphics display device (e.g. Xwindows). Graphics display in **BoA** is controlled by a software package called **BoGLi** (the **BoA** Graphic Library), which is described in Chapter 5. A few basic **BoGLi** commands which are needed in order to carry out the **BoA** plotting methods described in section 4.10 are thus described in this section.

### 4.9.1 Opening a plot window

Opening a graphic device is done with the `openDev()` command:

```
openDev()    # open a device, default: XWindow
open()       # alternatively, use one of the abbreviated commands
```

The default is to open an XWindow. You can use

```
open('?')
```

to get a list of all recognized devices. Alternatively, if you know which device you want you can enter it directly, for example

```
open('/ps')
```

You can also open a named PostScript file, here a colour PostScript file named *signal.ps*, with

```
open('signal.ps/CPS')
```

#### 4.9.2 Clearing a plot window

Clearing a plotting window is done with the `clear()` command:

```
clear()          # clear the active device
```

However, any plot command will first clear the active device before plotting a new graph, unless the *overplot=1* keyword is supplied.

#### 4.9.3 Closing a plot window

Closing a graphic device is done with the `closeDev()` command:

```
closeDev()      # open a device, default: XWindow
```

#### 4.9.4 Selecting an open device

METHOD: `selectDev`

DESCRIPTION: select an open device

### 4.10 Plotting and displaying data

#### 4.10.1 Displaying channel maps

METHOD: `chanMap(optional argument)`

DESCRIPTION: plot channel maps

OPTIONAL ARGUMENTS:

<i>chanList</i>	list of channels, of the form [1,2,3]
<i>flag</i>	flag values to consider
<i>oversamp</i>	oversampling factor (beam fwhm / pixel size). Default=2.
<i>center</i>	0/1 if set to 1 it will shift each map by the bolometer offset from the fits header. Thereby it shifts the source to the center of each channel map
<i>showRms</i>	compute and print rms/beam? (default: no)
<i>rmsKappa</i>	kappa in kappa-sigma clipping used to compute rms

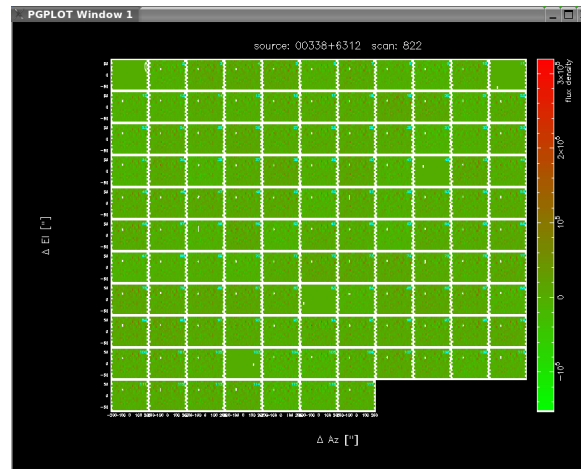


Figure 4.10.1: Default graphical outputs of a channel map of the source 00388+6312, including a wedge.

#### 4.10.2 Displaying/re-displaying a map

METHOD: `display(optional arguments)`

DESCRIPTION: display the reconstructed map in (Az,El)

OPTIONAL ARGUMENTS:

<i>weight</i>	plot the weight map instead of signal map
<i>coverage</i>	plot the rms map instead of signal map
<i>style</i>	the style used for the color (default idl4)
<i>caption</i>	the caption of the plot (default "")
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>limitsZ</i>	range of Z values to be plotted (comma separated values, in square brackets)
<i>wedge</i>	draw a wedge ? (default : yes)
<i>aspect</i>	keep the aspect ratio (default : yes)
<i>overplot</i>	should we overplot this image (default : no)
<i>doContour</i>	draw contour instead of map (default : no)
<i>levels</i>	the levels of the contours (default : intensity progression)
<i>labelContour</i>	label the contour (default : no)

#### Example:

Re-display a map after performing a smoothing (as in example in Section 3.5).  
`smooth(6./3600.) display(caption=data.ScanParam.caption())`

### 4.10.3 Plot the receiver parameters

METHOD: `plotArray` (*optional arguments*)

DESCRIPTION: plot the receiver layout

OPTIONAL ARGUMENTS:

<i>overplot</i>	overplot?
<i>num</i>	indicate channel numbers?

**Example:**

Plot the array layout with receiver numbers indicated. `plotArray (num=1)`

### 4.10.4 Plotting elevation versus azimuth

METHOD: `plotAzEl` (*optional arguments*)

DESCRIPTION: Plot elevation versus azimuth.

OPTIONAL ARGUMENTS:

<i>flag</i>	flag to be used (default = 0: all valid data; -1: plot all)
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	
<i>aspect</i>	

A more detailed description of the plotting related arguments can be found in [Section 5.5](#).

**Example:**

as for `plotAzimuth()`.

### 4.10.5 Plotting azimuth and elevation acceleration

METHOD: `plotAzElAcceleration` (*optional arguments*)

DESCRIPTION: Plot azimuth and elevation acceleration.

OPTIONAL ARGUMENTS:

*flag*        flag to be used (default = 0: all valid data; -1: plot all)  
*limitsX*    range of X values to be plotted (comma separated values, in square brackets)  
*limitsY*    range of Y values to be plotted (comma separated values, in square brackets)  
*style*       linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci*         colour index to be used (integer values)  
*overplot*  
*aspect*

A more detailed description of the plotting related arguments can be found in Section 5.5.

**Example:**

as for `plotAzimuth()`.

#### 4.10.6 Plotting elevation offset versus azimuth offset

METHOD: `plotAzElOffset` (*optional arguments*)

DESCRIPTION: Plot elevation offset versus azimuth offset.

OPTIONAL ARGUMENTS:

*flag*        flag to be used (default = 0: all valid data; -1: plot all)  
*limitsX*    range of X values to be plotted (comma separated values, in square brackets)  
*limitsY*    range of Y values to be plotted (comma separated values, in square brackets)  
*style*       linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci*         colour index to be used (integer values)  
*overplot*  
*aspect*

A more detailed description of the plotting related arguments can be found in Section 5.5.

**Example:**

as for `plotAzimuth()`.

#### 4.10.7 Plotting azimuth and elevation speed

METHOD: `plotAzElSpeed` (*optional arguments*)

DESCRIPTION: Plot azimuth and elevation speed.

OPTIONAL ARGUMENTS:

*flag* flag to be used (default = 0: all valid data; -1: plot all)  
*limitsX* range of X values to be plotted (comma separated values, in square brackets)  
*limitsY* range of Y values to be plotted (comma separated values, in square brackets)  
*style* linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci* colour index to be used (integer values)  
*overplot*  
*aspect*

A more detailed description of the plotting related arguments can be found in Section 5.5.

**Example:**

```
as for plotAzimuth().
```

#### 4.10.8 Plotting azimuth versus LST

METHOD: `plotAzimuth(optional arguments)`

DESCRIPTION: Plot the time series of the azimuth, i.e. azimuth versus LST.

OPTIONAL ARGUMENTS:

*flag* flag to be used (default = 0: all valid data; -1: plot all)  
*limitsX* range of X values to be plotted (comma separated values, in square brackets)  
*limitsY* range of Y values to be plotted (comma separated values, in square brackets)  
*style* linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci* colour index to be used (integer values)  
*overplot*  
*aspect*

A more detailed description of plotting related arguments can be found in Section 5.5.

**Example:**

```
azimuth(style='p', ci=2, limitsY=[-14,-13])
```

Plot azimuth versus LST (note the abbreviated form 'azimuth' used, see Table ??). Show individual plotted points (rather than lines), make plotted points red, and only plot azimuth (y axis) from -14 to -13 degrees.

#### 4.10.9 Plotting azimuth offset versus LST

METHOD: `plotAzimuthOffset(optional arguments)`

DESCRIPTION: Plot azimuth offset versus LST.

OPTIONAL ARGUMENTS:

<i>flag</i>	flag to be used (default = 0: all valid data; -1: plot all)
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	
<i>aspect</i>	

A more detailed description of the plotting related arguments can be found in Section 5.5.

**Example:**

```
as for plotAzimuth().
```

#### 4.10.10 Plot flux density of channels versus reference channel

METHOD: `plotCorrel` (*optional argument*)

DESCRIPTION: plot flux density of a list of channels vs. flux density of a reference channel

OPTIONAL ARGUMENTS:

<i>chanRef</i>	reference channel number (default : is the first in chanList)
<i>chanList</i>	list of channels, of the form [1,2,3]
<i>flag</i>	flag to be used
<i>skynoise</i>	plot against the skynoise of chanRef (default : no)
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	

#### 4.10.11 Plotting elevation versus LST

METHOD: `plotElevation` (*optional arguments*)

DESCRIPTION: Plot the time series of the elevation i.e. elevation versus LST.

OPTIONAL ARGUMENTS:

---

<i>flag</i>	flag to be used (default = 0: all valid data; -1: plot all)
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	
<i>aspect</i>	

A more detailed description of plotting related arguments can be found in Section 5.5.

**Example:**

```
as for plotAzimuth().
```

#### 4.10.12 Plotting elevation offset versus LST

METHOD: `plotElevationOffset` (*optional arguments*)

DESCRIPTION: Plot elevation offset versus LST.

OPTIONAL ARGUMENTS:

<i>flag</i>	flag to be used (default = 0: all valid data; -1: plot all)
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	
<i>aspect</i>	

A more detailed description of the plotting related arguments can be found in Section 5.5.

**Example:**

```
as for plotAzimuth().
```

#### 4.10.13 Plotting the FFT of the signal

METHOD: `plotFFT` (*optional arguments*)

DESCRIPTION: Plot a Fast Fourier Transform (FFT) of the signal

OPTIONAL ARGUMENTS:



<i>chanList</i>	list of channels, of the form [1,2,3]
<i>flag</i>	flag to be used
<i>labelX</i>	the X label; default is <i>Frequency [Hz]</i>
<i>labelY</i>	the Y label; default is <i>Amplitude (a.b.u/sqrt(Hz))</i>
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>plotphase</i>	plot phase instead of amplitude (default no)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	

**Example:**

Plot FFT for the first 9 channels.

```
plotFFT(range(10))
```

**4.10.14 Plot mean flux versus subscan number**

METHOD: `plotMean()` (*optional argument*)

DESCRIPTION: plot mean flux value vs. subscan number

OPTIONAL ARGUMENTS:

<i>chanList</i>	list of channels
<i>map</i>	plot as a 2D map?

**4.10.15 Plot mean channel values versus channel number**

METHOD: `plotMeanChan()` (*optional argument*)

DESCRIPTION: plot the MEAN value for each subscan against channel number

OPTIONAL ARGUMENTS:

<i>chanList</i>	list of channels
<i>flag</i>	flag to be used
<i>limitsX</i>	range of X values to be plotted (comma separated values, in square brackets)
<i>limitsY</i>	range of Y values to be plotted (comma separated values, in square brackets)
<i>style</i>	linestyle to be used ('p' or 'l', for points and solid line respectively)
<i>ci</i>	colour index to be used (integer values)
<i>overplot</i>	

**4.10.16 Plot flux Rms versus subscan number**

METHOD: `plotRms()` (*optional argument*)

DESCRIPTION: plot flux r.m.s. vs. subscan number

OPTIONAL ARGUMENTS:

*chanList* list of channels  
*map* plot as a 2D map?

#### 4.10.17 Plotting RMS versus channel number

METHOD: `plotRmsChan` (*optional arguments*)

DESCRIPTION: plot the RMS value for each subscan against channel number

OPTIONAL ARGUMENTS:

*chanList* list of channels  
*flag* flag to be used  
*limitsX* range of X values to be plotted (comma separated values, in square brackets)  
*limitsY* range of Y values to be plotted (comma separated values, in square brackets)  
*style* linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci* colour index to be used (integer values)  
*overplot*  
*subscan* if 0, plot rms of the complete scan (default); if 1, plot for each subscan and each channel

#### 4.10.18 Display start and end times of subscans

METHOD: `plotSubscan` ()

DESCRIPTION: generate a plot showing starting and ending times of subscans

#### 4.10.19 Plot subscans on the Az-El pattern

METHOD: `plotSubscanOffsets` ()

DESCRIPTION: Use four colours to show subscans on the Az, El pattern

#### 4.10.20 Show a reconstructed map in (Az,El) or (Ra,Dec)

#### 4.10.21 Plotting flux density versus LST

METHOD: `signal` (*optional argument*)

DESCRIPTION: Plot the time series of the flux density i.e. flux density versus LST.

OPTIONAL ARGUMENTS:

*chanList* list of channels, of the form [1,2,3]  
*flag* flag to be used  
*mjd* if set, use mjd instead of lst  
*limitsX* range of X values to be plotted (comma separated values, in square brackets)  
*limitsY* range of Y values to be plotted (comma separated values, in square brackets)  
*style* linestyle to be used ('p' or 'l', for points and solid line respectively)  
*ci* colour index to be used (integer values)  
*overplot*

A more detailed description of the plotting related arguments can be found in [Section 5.5](#).

**Example:**

```

signal(chanList=[18,19,20], mjd=1, style='p', ci=2)
signal([18,19,20], mjd=1, style='p', ci=2)

```

## 4.11 Data handling

### 4.11.1 Get a list of valid channels

METHOD: `checkChanList` (*optional argument*)

DESCRIPTION: Return a list of valid channels

OPTIONAL ARGUMENTS:

*inList* list of channel numbers to get, or empty list to get the complete list  
of unflagged channels, or 'all' or 'al' or 'a' to get the complete list of channels  
*flag* flag value to be selected (default = 0)

### 4.11.2 Get pixel values

METHOD: `getPixel()`

DESCRIPTION: get pixel values using mouse

Click left to get one pixel, mid to get average over 9, right to exit (on Data array only).

### 4.11.3 Print the current list of channels

METHOD: `printChannels()`

DESCRIPTION: Print the current list of channels

#### 4.11.4 Selecting channels

METHOD: `setCurrChanList (optional argument)`

DESCRIPTION: Set a channel or a list of channels to be treated.

OPTIONAL ARGUMENTS:

<i>chanList:</i>	list of channel numbers, of the form: [1,2,3]
'all'... 'al'...'a'	set current list to all possible channels
'?'	get current list of channels (default)

#### Example:

```
Using the abbreviated form channels() (see Table 4.1): channels([1,2,3])
channels(chanList=[1,2,3])
channels('all')
channels('?')
```

## 4.12 MB-Fits to FITS file conversion

To convert an MB-Fits file to a FITS file in the MAMBO format you can use the command `mambo()`. The current version does NOT use the data contained in the data object in **BoA**, but reads the input file (with default name = `BoaB.currData.FileName`) and converts it to the Mambo format. Therefore, this procedure is somewhat decoupled from **BoA**.

## 4.13 User methods for selecting files and directories

### 4.13.1 Listing the contents of the input directory

METHOD: `listInDir()`

DESCRIPTION: list the contents of the input directory

### 4.13.2 Resetting the CurrentList

METHOD: `resetCurrentList()`

DESCRIPTION: reset the CurrentList to the complete List

### 4.13.3 Setting the input directory

METHOD: `setInDir()`

DESCRIPTION: set the input directory

**Example:**

```
setInDir('inputDirectory')
```

**4.13.4 Setting the output directory**

METHOD: `setOutDir()`

DESCRIPTION: set the input directory

**Example:**

```
setOutDir('outputDirectory')
```

**4.13.5 Setting the project ID**

METHOD: `setProjectID()`

DESCRIPTION: set the current project ID

**Example:**

```
setProjectID('projectID')
```

**4.14 Miscellaneous methods****4.14.1 Updating offset and gain values from a file**

METHOD: `updateRCP(rcpFile)`

DESCRIPTION: update only offsets and gains from the content of a file

INPUT:

*rcpFile*    complete name of file to read in

**4.15 Scripts**

As **BoA** provides the full functionality of Python this allows the use of scripts. Scripts can be run with the `execfile()` function where the name of the file has to be given as string argument. The suffix of the file is arbitrary.

**Example:**

If you want to have a look at the time series of channels 10 to 30 successively, create the following script with your preferred editor. Note that in Python the contents of the for loop (like if blocks, method definitions, etc.) have to be indented.

```
# testBoa.py
indir('../Fits/')      # set the input directory
read('test')           # read file test.fits
op()                   # open graphic display
for i in range(10,31): # start a for loop, the indentation in
                        # the following lines is mandatory
    sig([i])           # plot time series
    raw_input()        # wait for <Return>
```

To run the script type:

```
execfile('testBoa.py')
```

#### 4.15.1 Example scripts

In order to demonstrate some of the basic functionalities of **BoA** three demonstration scripts are provided: *ExampleMap.py*, *ExamplePointing.py* and *ExampleFocus.py*. These can be found in the directory `/home/user/boa/examples/` and are described in detail in Chapter 3. Run the scripts by typing:

```
execfile('/home/user/boa/examples/ExampleMap.py')
```

## 4.16 Commands in alphabetical order

<b>chanMap</b>	plot channel maps
<b>checkChanList</b>	return a list of valid channels
<b>clear</b>	clear the active plot window
<b>closeDev</b>	close one device
<b>computeRms</b>	compute rms/beam in a map
<b>computeWeight</b>	compute weights (default is use $1/\text{rms}^2$ )
<b>despike</b>	Flag yet unflagged data below 'below'*rms and above 'above'*rms
<b>display</b>	show the reconstructed maps in (Az,El)
<b>doMap</b>	construct a map in (Az,El) coordinates
<b>flagChannels</b>	flag a list of channels
<b>flagMJD</b>	flag data by time interval
<b>flagPosition</b>	flag a position in the sky within a given radius
<b>flagRms</b>	Flag channels with rms below or above respective given values
<b>flagSubscan</b>	flag certain subscans
<b>getPixel</b>	get pixel values using mouse
<b>horizontalMap</b>	construct a map in (Az,El) coordinates
<b>listInDir</b>	list the input directory
<b>medianBaseline</b>	baseline: Remove median value per channel and per sub-scan
<b>medianNoiseRemoval</b>	remove median noise from the data
<b>openDev</b>	open a graphic device
<b>plotArray</b>	plot the receiver layout
<b>plotAzEl</b>	plot elevation versus azimuth
<b>plotAzElAcceleration</b>	plot azimuth and elevation acceleration
<b>plotAzElOffset</b>	plot elevation offset versus azimuth offset
<b>plotAzElSpeed</b>	plot azimuth and elevation speed
<b>plotAzimuth</b>	plot azimuth versus LST
<b>plotAzimuthOffset</b>	plot azimuth offset versus LST
<b>plotCorrel</b>	plot signal vs. reference channel
<b>plotElevation</b>	plot elevation versus LST
<b>plotElevationOffset</b>	plot elevation offset versus LST
<b>plotFFT</b>	plot FFT of signal
<b>plotMean</b>	plot mean flux values vs. subscan numbers
<b>plotMeanChan</b>	plot mean value for each subscan vs. chan. number
<b>plotRms</b>	plot rms flux values vs. subscan numbers

---

<b>plotRmsChan</b>	plot rms value for each subscan vs. chan. number
<b>plotSubscan</b>	generate a plot showing starting and ending times of sub-scans
<b>plotSubscanOffsets</b>	Use four colours to show subscans on the Az, El pattern
<b>polynomialBaseline</b>	remove a polynomial baseline from the data
<b>printCurrChanList</b>	print the current channel list
<b>read</b>	read in a file
<b>resetCurrentList</b>	reset the CurrentList to the complete list
<b>saveMambo</b>	convert MB-Fits file to MAMBO format
<b>selectDev</b>	select an open device
<b>setCurrChanList</b>	select list of channels
<b>setInDir</b>	set the input directory
<b>setOutDir</b>	set the output directory
<b>setProjectID</b>	set the project ID
<b>showMap</b>	show the reconstructed map in (Az,El) or (Ra,Dec)
<b>signal</b>	plot the time series of the data (flux density versus LST)
<b>smoothBy</b>	smooth the image with a 2D gaussian of gived FWHM
<b>solveFocus</b>	compute the optimal focus position
<b>solvePointingOnMap</b>	compute the offset on the data.Map object
<b>statistics</b>	prints the statistics
<b>unflag</b>	unflag data
<b>unflagChannels</b>	unflag a list of channels
<b>updateRCP</b>	update offsets and gains from the content of a file



## 4.17 Commands in functional order

### 4.17.1 Plotting and displaying

<b>chanMap</b>	plot channel maps
<b>display</b>	show the reconstructed maps in (Az,El)
<b>plotArray</b>	plot the receiver layout
<b>plotAzEl</b>	plot elevation versus azimuth
<b>plotAzElAcceleration</b>	plot azimuth and elevation acceleration
<b>plotAzElOffset</b>	plot elevation offset versus azimuth offset
<b>plotAzElSpeed</b>	plot azimuth and elevation speed
<b>plotAzimuth</b>	plot azimuth versus LST
<b>plotAzimuthOffset</b>	plot azimuth offset versus LST
<b>plotCorrel</b>	plot signal vs. reference channel
<b>plotElevation</b>	plot elevation versus LST
<b>plotElevationOffset</b>	plot elevation offset versus LST
<b>plotFFT</b>	plot FFT of signal
<b>plotMean</b>	plot mean flux values vs. subscan numbers
<b>plotMeanChan</b>	plot mean value for each subscan vs. chan. number
<b>plotRms</b>	plot rms flux values vs. subscan numbers
<b>plotRmsChan</b>	plot rms value for each subscan vs. chan. number
<b>plotSubscan</b>	generate a plot showing starting and ending times of sub-scans
<b>plotSubscanOffsets</b>	Use four colours to show subscans on the Az, El pattern
<b>showMap</b>	show the reconstructed map in (Az,El) or (Ra,Dec)
<b>signal</b>	plot the time series of the data (flux density versus LST)

### 4.17.2 Device handling

<b>clear</b>	clear the active plot window
<b>closeDev</b>	close one device
<b>openDev</b>	open a graphic device
<b>selectDev</b>	select an open device

### 4.17.3 Pointing and focus

<b>solveFocus</b>	compute the optimal focus position
<b>solvePointingOnMap</b>	compute the offset on the data.Map object

**4.17.4 Flagging and despiking data**

<b>despike</b>	flag yet unflagged data below 'below'*rms and above 'above'*rms
<b>flagChannels</b>	flag a list of channels
<b>flagMJD</b>	flag data by time interval
<b>flagPosition</b>	flag a position in the sky within a given radius
<b>flagRms</b>	flag channels with rms below or above respective given values
<b>flagSubscan</b>	flag certain subscans
<b>unflag</b>	unflag data
<b>unflagChannels</b>	unflag a list of channels

**4.17.5 Map making**

<b>doMap</b>	construct a map in (Az,El) coordinates
<b>horizontalMap</b>	construct a map in (Az,El) coordinates

**4.17.6 Baseline subtraction, sky removal and statistics**

<b>computeRms</b>	compute rms/beam in a map
<b>computeWeight</b>	compute weights (default is use $1/\text{rms}^2$ )
<b>medianBaseline</b>	baseline: Remove median value per channel and per subscan
<b>medianNoiseRemoval</b>	remove median noise from the data
<b>polynomialBaseline</b>	remove a polynomial baseline from the data
<b>smoothBy</b>	smooth the image with a 2D gaussian of given FWHM
<b>statistics</b>	prints the statistics

**4.17.7 File handling**

<b>read</b>	read in a file
<b>saveMambo</b>	convert MB-Fits file to MAMBO format

**4.17.8 Data handling**

<b>checkChanList</b>	return a list of valid channels
<b>getPixel</b>	allow user to get pixel values using mouse
<b>printCurrChanList</b>	print the current channel list
<b>setCurrChanList</b>	select list of channels

**4.17.9 Selecting files and directories**

<b>listInDir</b>	list the input directory
<b>resetCurrentList</b>	reset the CurrentList to the complete list
<b>setInDir</b>	set the input directory
<b>setOutDir</b>	set the output directory
<b>setProjectID</b>	set the project ID

**4.17.10 Misc.**

<b>updateRCP</b>	update offsets and gains from the content of a file
------------------	---

## 4.18 Abbreviations

As we have noted already, user methods are abbreviations of the full methods. For example, the method `DeviceHandler.openDev()` can be called by the user method `open()`. For further convenience, most user methods can also be called by even shorter abbreviations of the user methods (in this example `op()` is all that is needed). A list of user methods and their abbreviations is given in Table 4.1.

Command	Abbreviations
chanMap	ChanMap - chanmap
checkChanList	checkChannels - checkChan
clear	cle - cl
closeDev	close - clo - cls
computeWeight	computeweight - weight
despike	dspike
display	mapdisp - mapdisplay
doMap	domap
findInDir	find - fd
flag	
flagChannels	flagCh - flagC - fCh
flagLon	
flagMJD	
flagPosition	flagPos
flagRms	
flagSubscan	flagSub
getPixel	getPix
listInDir	ils - inls
mapping	horizontalMap - hrMap
maprms	computeRms
medianBaseline	medianBase - medianbase
medianNoiseRemoval	mediannoise
openDev	open - op
plotArray	plotarray
plotAzEl	azel
plotAzElAcceleration	azelaccel - azelac
plotAzElOffset	azeloff - azelo
plotAzElSpeed	azelspeed - azelsp
plotAzimuth	azimuth - azimuth - az
plotAzimuthOffset	azimuthOffset - azimuthoff - azo
plotCorrel	plotcorrel - plotcor - plotCor
plotElevation	elevation - elev - el
plotElevationOffset	elevationOffset - eleoff - elo

Table 4.1: List of user methods with abbreviations. Don't forget to add the round brackets () at the end of the commands.

Command	Abbreviations
plotFFT	
plotMean	plotmean
plotMeanChan	plotmeanchan
plotRms	plotrms
plotRmsChan	plotrmschan
plotSubscan	plotSub
plotSubscanOffsets	plotSubOff
polynomialBaseline	baseline - base
printCurrChanList	printChannels - printChan
read	
resetCurrentList	resetCurrList - rls
selectDev	device - dev
selectInDir	select - slt
setCurrChanList	channels - channel - chan
setInDir	indir - ind
setInFile	infile - inf
setOutFile	outfile - outf
setOutDir	outdir - outd
setProjectID	setproj - proj
showMap	
signal	signa - sign - sig
solveFocus	solvefocus - solveFoc - solvefoc
solvePointingOnMap	solvepointing - solvepoint - solvepoin - solvepoi
smoothBy	smooth
statistics	stats - stat
unflag	
unflagChannels	unflagCh - unflagC - ufCh
updateRCP	

Table 4.1: *continued*

---

## 5. BoGLi: THE BoAGRAPHIC LIBRARY

---

### 5.1 Introduction

The **BoA** Graphic Library (**BoGLi**) is an object-oriented software package for the graphical display of data. It is written in Python and uses **pgplot**, the python binding to **pgplot**. The main parts (classes) of the software are self-consistent and may independently be used from any python programme. Nevertheless, **BoGLi** comes with features which especially customise its use for the display of astronomical data from multi-channel receivers. Its main goal is to provide a graphic tool tailored for the use with **BoA** for the display of data from LaBoCa, Simba and Mambo.

### 5.2 Command handling

**BoGLi** has its own command handler. Nevertheless, anytime the **BoA** command handler encounters a graphic command this is automatically passed to the **BoGLi** command handler. Therefore, the user does not have to care about the separation between **BoA** and **BoGLi** commands. Table 5.1 gives an overview of some of the available commands.

**BoGLi** provides a variety of attributes that may be changed by the user. The attribute name is then used as command followed by the desired value as argument (see Sect. ?? for details.)

Table 5.1: List of useful **BoGLi** commands.

<code>DeviceHandler.openDev</code>	open a device
<code>DeviceHandler.closeDev</code>	close a device
<code>Plot.clear</code>	clear the active plot window
<code>DeviceHandler.selectDev</code>	select a device
<code>DeviceHandler.resizeDev</code>	resize the plotting area, after plot window resized using mouse
<code>Plot.plot</code>	make a single plot
<code>MultiPlot.plot</code>	plot multiple plots
<code>Plot.draw</code>	draw on an image
<code>MultiPlot.draw</code>	draw on plots of multiple channels

### 5.3 Device handling

BoGLi is based on `pgplot` and as a consequence the number and type of available devices depends on the actual configuration. A list of supported devices is given at <http://www.astro.caltech.edu/~tjp/pgplot/devices.html>. During installation the device drivers have to be selected by editing the file *drivers.list*. As many device drivers are available on selected operating systems only, you should ensure that drivers you do not want are commented out (place ! in column 1) to avoid installation failures. A version of *drivers.list* used for a Linux PC can be found in Sect ??.

The command handler of **BoGLi** provides a set of commands to manage output devices. A detailed description of these commands is given below.

#### 5.3.1 Opening a plot window

**DESCRIPTION:** Open a graphics device for `pgplot` output and make it the current device. The default, when no argument is provided, is to open an XWindow.

**USAGE:** `DeviceHandler.openDev (optional argument)`

The relevant abbreviations can also be used (see Table 4.1).

**OPTIONAL ARGUMENT:** *pgplot device type*

If the device is opened successfully, it becomes the selected device to which graphics output is directed until another device is selected (see 5.3.4) or the device is closed (see 5.3.2). If no device argument is specified PGPLOT will open the default graphics device (an XWINDOW). Alternatively, the graphics device may be selected using any of the following as arguments:

- (1) A complete device specification of the form 'device/type' or 'file/type', where /type is one of the allowed PGPLOT device types (installation-dependent, e.g. /xwindow) and 'device' or 'file' is the name of a graphics device or disk file appropriate for this type. The 'device' or 'file' may contain '/' characters; the final '/' delimits the 'type'. If necessary to avoid ambiguity, the 'device' part of the string may be enclosed in double quotation marks.

**Example:** `'plot.ps/ps', 'dir/plot.ps/ps', '"dir/plot.ps"/ps', 'user:[tjp.plots]plot.ps/PS'`

- (2) A device specification of the form '/type', where /type is one of the allowed PGPLOT device types, e.g. /xwindow. PGPLOT supplies a default file or device name appropriate for this device type.

**Example:** `'/ps'` (PGPLOT interprets this as 'pgplot.ps/ps')

- (3) A device specification with '/type' omitted; in this case the type is taken from the environment variable PGPLOT\_TYPE, if defined (e.g., `setenv PGPLOT_TYPE PS`). Because of possible confusion with '/' in file-names, omitting the device type in this way is not recommended.

**Example:** `'plot.ps'` (if PGPLOT\_TYPE is defined as 'ps', PGPLOT interprets this as 'plot.ps/ps')

- (4) A blank string (' '); in this case, PGOPEN will use the value of environment variable PGPLOT\_DEV as the device specification, or '/NULL' if the environment variable is undefined.

**Example:** ' ' (if PGPLOT\_DEV is defined)

- (5) A single question mark, with optional trailing spaces, i.e. ('?'). In this case, PGPLOT will prompt the user to supply the device specification, with a prompt string of the form 'Graphics device/type (? to see list, default XXX):' where 'XXX' is the default (value of environment variable PGPLOT\_DEV).

**Example:** ' ? '

- (6) A non-blank string in which the first character is a question mark (e.g. '?Device: '); in this case, PGPLOT will prompt the user to supply the device specification, using the supplied string as the prompt (without the leading question mark but including any trailing spaces).

**Example:** '?Device specification for PGPLOT: '

In cases (5) and (6), the device specification is read from the standard input. The user should respond to the prompt with a device specification of the form (1), (2), or (3). If the user types a question-mark in response to the prompt, a list of available device types is displayed and the prompt is re-issued. If the user supplies an invalid device specification, the prompt is re-issued. If the user responds with an end-of-file character, e.g., ctrl-D in UNIX, program execution is aborted; this avoids the possibility of an infinite prompting loop. A programmer should avoid use of PGPLOT-prompting if this behavior is not desirable.

The device type is case-insensitive (e.g., '/ps' and '/PS' are equivalent). The device or file name may be case-sensitive in some operating systems.

### 5.3.2 Closing a plot window

**DESCRIPTION:** Close a plotting device. The default, where no argument is supplied, is to close the current device.

**USAGE:** `DeviceHandler.closeDev (optional argument)`

**OPTIONAL ARGUMENT:**

*device number* (integer)  
'all'  
'current'...'curre'...'cur'

**Example:**

<code>DeviceHandler.closeDev(2)</code>	Close the device with identifier 2
<code>DeviceHandler.closeDev('all')</code>	close all devices
<code>DeviceHandler.closeDev('current')</code>	close current device (the default if no argument specified)

### 5.3.3 Clearing a plot window

**DESCRIPTION:** Clear the output of the current device. To clear the output of a different device change to that device first (see [5.3.4](#)).



USAGE: `Plot.clear()`

### 5.3.4 Selecting a device

DESCRIPTION: Select an open device for graphical output. The selected device has to be previously opened with *open* (see 5.3.1).

USAGE: `DeviceHandler.selectDev(argument)`

ARGUMENT: *device number* (integer)

#### Example:

```
DeviceHandler.selectDev(2)    Make device number 2 the current device for graphical output
```

### 5.3.5 Resizing a device

DESCRIPTION: Resize the plotting area after resizing of the graphics display window using the mouse. This is applicable to some interactive devices (e.g. /xwindow).

USAGE: `DeviceHandler.resizeDev()`

## 5.4 Plotting graphics

This section lists some of the graphics plotting capabilities of **BoGLi**.

### 5.4.1 Plotting single plots

DESCRIPTION: Make a single plot of x versus (optional) y.

USAGE: `Plot.plot( dataX, [ dataY, limitsX, limitsY, labelX, labelY, caption, style, ci, width, overplot, aspect, logX, logY, noData ] )`

ARGUMENTS:

*dataX* values to plot along X

*dataY* values to plot along Y (optional - default: plot dataX vs. running number)

OPTIONAL ARGUMENTS:

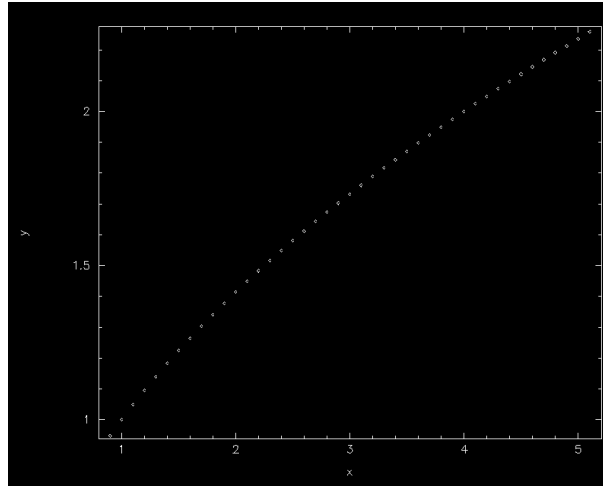


Figure 5.4.1: Example 1 of graphics produced using Plot.plot

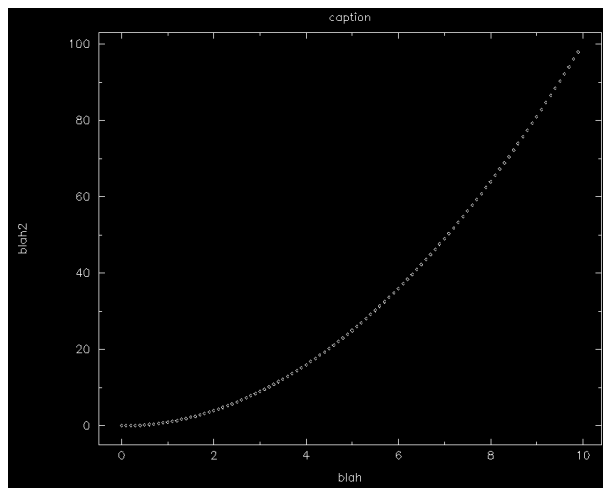


Figure 5.4.2: Example 2 of graphics produced using Plot.plot

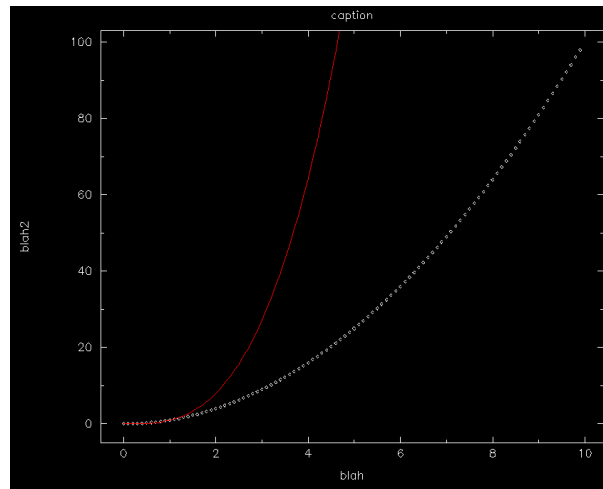


Figure 5.4.3: Example 3 of graphics produced using Plot.plot

<i>limitsX</i>	limits to use in X for the plot
<i>limitsY</i>	limits to use in Y for the plot
<i>labelX</i>	x label (default 'x')
<i>labelY</i>	y label (default 'y')
<i>caption</i>	the caption of the plot (default '')
<i>style</i>	the style used for the plot ('l': line, 'p': point (default), 'b': histogram)
<i>ci</i>	color index (default 1)
<i>width</i>	linewidth (default 0 = use previous)
<i>aspect</i>	keep the aspect ratio in 'physical' unit
<i>overplot</i>	set overplot=1 to overplot (default no)
<i>logX</i>	set logX=1 to use a log scale (default no)
<i>logY</i>	set logY=1 to use a log scale (default no)

These are also described in Section ???. Note *dataY* is also optional – if no *dataY* is supplied the default is to plot *dataX* versus running number.

#### Example:

```
x = Numeric.array(range(100), Numeric.Float)/10
```

```
Plot.plot(x, Numeric.sqrt(x), limitsX=[1, 5])
```

Note that Y limits are then computed according to this X range.

The graphic output produced in this case is shown in Figure 5.4.1.

#### Example:

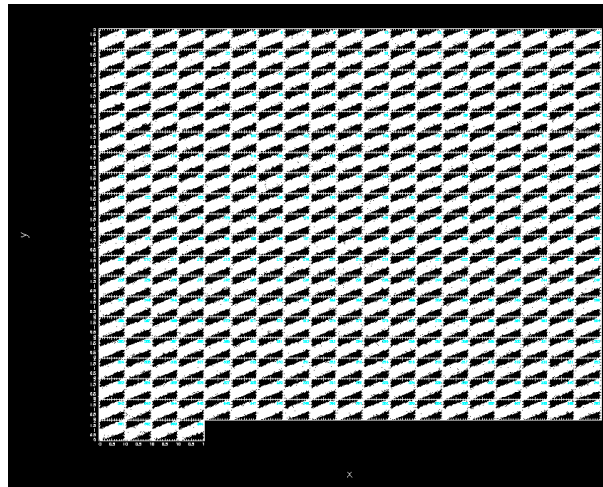


Figure 5.4.4: Example of graphics produced using MultiPlot.plot

```
Plot.plot(x, x*x, labelX='blah', labelY='blah2', caption='caption')
```

Note that plot clear the screen first, you need to use the new 'overplot' keyword (see below).

The graphic output produced in this case is shown in Figure 5.4.2.

#### Example:

```
Plot.plot(x, x*x*x, overplot=1, ci=2, style='l')
```

The graphic output produced in this case is shown in Figure 5.4.3.

### 5.4.2 Plotting multiple channels

DESCRIPTION: Make a plot of x versus (optional) y for several channels simultaneously.

USAGE: `MultiPlot.plot(chanList, dataX, dataY, [ limitsX, limitsY, labelX, labelY, caption, style, ci, overplot, logX, logY, nan ])`

ARGUMENTS:

<i>chanList</i>	list of channels, of the form [1,2,3]
<i>dataX</i>	values to plot along X
<i>dataY</i>	values to plot along Y

OPTIONAL ARGUMENTS:

<i>limitsX</i>	limits to use in X for the plot
<i>limitsY</i>	limits to use in Y for the plot
<i>labelX</i>	x label (default 'x')
<i>labelY</i>	y label (default 'y')
<i>caption</i>	the caption of the plot (default '')
<i>style</i>	the style used for the plot ('l': line, 'p': point (default), 'b': histogram)
<i>ci</i>	color index (default 1)
<i>overplot</i>	set overplot=1 to overplot (default no)
<i>logX</i>	set logX=1 to use a log scale (default no)
<i>logY</i>	set logY=1 to use a log scale (default no)

These are also described in Section ??.

#### Example:

```
n_point = 365
chanlist=range(n_point)

x2 = RandomArray.random([n_point,n_point])
y2 = RandomArray.random([n_point,n_point])

MultiPlot.plot(chanlist,x2,y2+x2,style='p')
```

The graphic output produced in this case is shown in Figure [5.4.4](#).

### 5.4.3 Drawing on an image

DESCRIPTION: Draw on an image

USAGE: `Plot.draw( map_array, [ sizeX, sizeY, WCS, limitsX, limitsY, limitsZ, nan, labelX, labelY, caption, style, contrast, brightness, wedge, overplot, aspect, doContour, levels, labelContour ] )`

ARGUMENTS:

*map\_array* map to display

OPTIONAL ARGUMENTS:

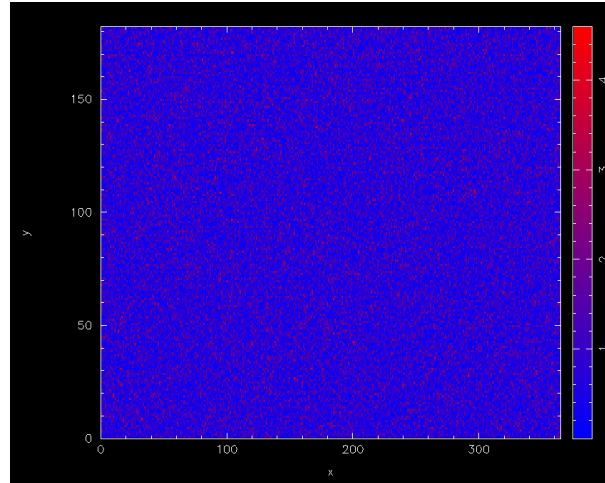


Figure 5.4.5: Example 1 of graphics produced using Plot.draw

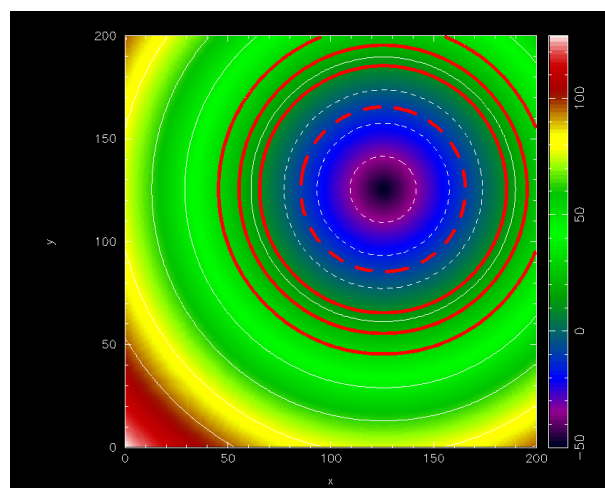


Figure 5.4.6: Example 2 of graphics produced using Plot.draw: drawing contours

<i>sizeX</i>	the 'physical' size of the array (default pixel numbers), defined by the center of the two extreme pixels
<i>sizeY</i>	the 'physical' size of the array (default pixel numbers), defined by the center of the two extreme pixels
<i>limitsX</i>	limits to use in X for the plot
<i>limitsY</i>	limits to use in Y for the plot
<i>nan</i>	set =1 if NaN are present in the array
<i>labelX</i>	x label (default 'x')
<i>labelY</i>	y label (default 'y')
<i>caption</i>	the caption of the plot (default '')
<i>style</i>	the color used for the plot (default 'g2r', see <code>Plot.Plot.setImaCol()</code> )
<i>wedge</i>	set <code>wedge=1</code> to draw a wedge (default no)
<i>aspect</i>	keep the aspect ratio in 'physical' unit
<i>overplot</i>	set <code>overplot=1</code> to overplot (default no)
<i>doContour</i>	set =1 to draw contour instead of map (default no)
<i>levels</i>	the levels for the contours (default <code>nContour</code> , within <code>plotLimitsZ</code> )
<i>labelContour</i>	set =1 to label the contours (default no)

These arguments are also described in Section ??.

#### Example:

```
n_point = 365

mapping = Numeric.absolute(RandomArray.standard_normal([n_point,n_point/2]))

Plot.draw(mapping, style='b2r', wedge=1)

# You can also define 'physical' unit for your plot and still use
# limitsX/Y and aspect:

Plot.draw(mapping, sizeX=[-1,1], sizeY=[-2,2], limitsY=[-1,1], aspect=1, wedge=1)
```

The graphic output produced in this case is shown in Figure 5.4.5.

#### Example:

You can also use `Plot.draw()` to plot contours.

```
def dist(x,y):
    return (x-125)**2+(y-125)**2
```

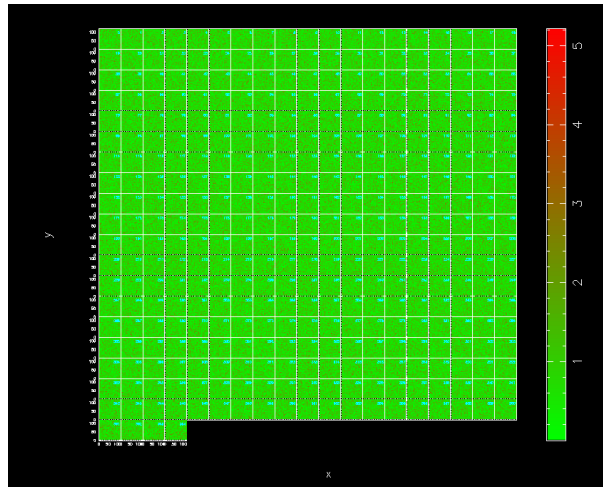


Figure 5.4.7: Example of graphics produced using MultiPlot.draw

```

image = Numeric.sqrt(Numeric.fromfunction(dist, (200,200)))-50

Plot.draw(image,wedge=1,aspect=1,style='rainbow')      # display an image
Plot.draw(image,doContour=1,overplot=1)               # overlay some contours
Plot.contour['color'] = 2                             # change the colour and
Plot.contour['linewidth'] = 10                       # linewidth attributes

Plot.draw(image,doContour=1,overplot=1,levels=[-10,10,20,30]) # plot some
# more contours with the new attributes

```

The graphic output produced in this case is shown in Figure 5.4.6.

#### 5.4.4 Drawing on plots of multiple channels

DESCRIPTION: Draw on a multi-channel image

USAGE: `MultiPlot.plot.draw( chanList, map_arrays, [ sizeX, sizeY, WCS, limitsX, limitsY, limitsZ, nan, labelX, labelY, caption, style, contrast, brightness, wedge, overplot ] )`

ARGUMENTS:

*chanList*      list of channels  
*map\_arrays*    lits of map to display

OPTIONAL ARGUMENTS:



---

<i>sizeX</i>	the 'physical' size of the array (default pixel numbers)
<i>sizeY</i>	the 'physical' size of the array (default pixel numbers)
<i>limitsX</i>	limits to use in X for the plot
<i>limitsY</i>	limits to use in Y for the plot
<i>labelX</i>	x label (default 'x')
<i>labelY</i>	y label (default 'y')
<i>caption</i>	the caption of the plot (default '')
<i>style</i>	the color used for the plot (default 'g2r', see <code>Plot.Plot.setImaCol()</code> )
<i>wedge</i>	set <code>wedge=1</code> to draw a wedge (default no)
<i>overplot</i>	set <code>overplot=1</code> to overplot (default no)

These are also described in Section ??.

**Example:**

```
mapping_array = []
n_map = 365
for i in range(n_map):
    mapping_array.append(Numeric.absolute(RandomArray.standard_normal([120,
MultiPlot.draw(range(n_map), mapping_array, wedge=1)
```

The graphic output produced in this case is shown in Figure [5.4.7](#).

## 5.5 Keywords

**BoGLi** provides a variety of parameters which allow the graphical output to be customised, as regards primitives such as colours, linestyles, character sizes, as well as text output and general appearance.

**ci**      *colour index*

The colour index is an integer in the range 0 to a device-dependent maximum. The default colour index is 1, usually white on a black background for monitor displays or black on a white background for printed hardcopies. Colour index 0 corresponds to the background colour. If the requested color index is not available on the selected device, colour index 1 will be used.

**ls**      *line style*

The line style is an integer in the range 1 to 5 with the following codes:

- 1: full line
- 2: dashed
- 3: dot-dash-dot-dash
- 4: dotted
- 5: dash-dot-dot-dot

The line style does not affect graph markers, text, or area fill.

**lw**      *line width*

The line width is specified in units of 1/200 (0.005) inch (about 0.13 mm) and must be an integer in the range 1-201. This parameter affects lines, graph markers and text.

**limitsX**    *limits to use in X for the plot*

**limitsY**    *limits to use in Y for the plot*

**labelX**    *x label*  
(default 'x')

**labelY**    *y label (default 'y')*

**caption**    *caption label*  
(default ' ')

**style**      *linestyle*  
(**'l'**: line, **'p'**: point (default), **'b'**: histogram)

**width**      *linewidth*  
(default 0 = use previous)

---

<b>aspect</b>	<i>aspect ratio</i> keep the aspect ratio in 'physical' unit
<b>overplot</b>	<i>allow/prohibit overplotting</i> set overplot=1 to overplot (default no)
<b>logX</b>	<i>logarithmic scale</i> set logX=1 to use a log scale (default no)
<b>logY</b>	<i>logarithmic scale</i> set logY=1 to use a log scale (default no)
<b>sizeX</b>	<i>set the 'physical' size of the array</i> the 'physical' size of the array (default pixel numbers), defined by the center of the two extreme pixels
<b>sizeY</b>	<i>set the 'physical' size of the array</i> the 'physical' size of the array (default pixel numbers), defined by the center of the two extreme pixels
<b>nan</b>	set =1 if NaN are present in the array
<b>wedge</b>	set wedge=1 to draw a wedge (default no)
<b>doContour</b>	<i>draw contours</i> set =1 to draw contour instead of map (default no)
<b>levels</b>	<i>set the levels for the contours</i> the levels for the contours (default nContour, within plotLimitsZ)
<b>labelContour</b>	<i>label the contours</i> set =1 to label the contours (default no)

**Part II**

**Reference Manual**

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## 6. DATA ORGANISATION

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### 6.1 Data input: the MB-FITS format

A complete description of the Multi-Beam FITS Raw Data Format is given in the reference document APEX-MPI-IFD-0002. In this section, we only give a brief description of this file format.

#### 6.1.1 The hierarchy for a full scan

For a given observing sequence, corresponding to one scan, a set of tables are generated and stored in a hierarchical way in the MB-FITS format. Three tables are created on top of this hierarchy, where informations related to the full scan are gathered:

- Primary header: here, some general informations are stored, such as telescope name, project ID, date of observation start, versions of MB-FITS format and FitsWriter software
- SCAN-MBFITS: the header of this table contains a description of the scan pattern (type, geometry, line length in case of a raster map...), the source name and coordinates, together with a description of the referential used, and some generic informations about the telescope (coordinates, pointing coefficients). In addition, a binary table lists the names of frontend-backend (hereafter FEBE) combinations in use for this observation.
- FEBEPAR-MBFITS: one such table is created for each FEBE in use (in general, only one FEBE is active for bolometer observing). It contains the FEBE name and the number of available channels for this FEBE in its header. The associated binary table gives all relevant information about the instrument: relative gains, positions, gain/attenuation factors, polarisation angles...

#### 6.1.2 Tables for each subscan

For each subscan within a scan, three tables are generated:

- MONITOR-MBFITS: this table gathers all the monitoring information sent by the control system during the observation. Each datapoint has an associated timestamp in MJD. In particular, this monitor stream contains commanded and actual telescope positions sampled every 48 ms. It also contains data related to the weather conditions, the subreflector angle and position, and the LST values.
- DATAPAR-MBFITS: this table also contains the telescope positions, subreflector angles and positions, and LST values, but interpolated to the timestamps corresponding to the data stream.

It also contains a PHASE column, which can for example contains a succession of “ON” and “OFF” for a wobbler-switching observation.

- **ARRAYDATA-MBFITS**: here the raw data are stored. While some basic informations are stored in the header (e.g. central frequency of the observation), the binary table only contains two columns: the timestamps (in MJD), and a vector with length equal to the number of channels in use containing the raw data for each integration.

**Note:** in case several FEBE are in use at the same time, then a DATAPAR table and an ARRAYDATA table are generated for each subscan and for each FEBE.

## 6.2 BoAData objects

The manipulation of data within BoA is done with data objects of one class that inherits from the DataEntity class (Sect. 4.1.2; see also Section ??). Such objects contain the current version of the data, as well as associated parameters related to the scan and to the bolometer array. On top of this, the DataAna and Map classes define additional attributes, as described in the next subsections.

### 6.2.1 DataEntity

A DataEntity object has a number of attributes, listed in the following tables. Two of them are objects of classes BolometerArray and ScanParameter.

#### BolometerArray

The BolometerArray object defines the attributes listed in Table 6.1. They are read in from the file, or computed when reading, except for CurrChanList (contains the current list of channels on which any processing or plotting function is applied) and Flags (can be altered by the user).

#### Telescope

Attributes of a Telescope object are shown in Table 6.2.

#### ScanParam

Attributes of the ScanParam object (class ScanParameter) are listed in Table 6.3.

#### Data arrays

In addition to the scan parameters and bolometer array related informations, a DataEntity object contains some general informations about the observation, and 2D arrays of data and related numbers, with sizes number of pixels in use  $\times$  number of integrations. These are described in Table 6.4.

Table 6.1: Attributes of a BolometerArray object

Name	Type	Description
Telescope	object	see Table 6.2
FeBe	string	Frontend-Backend name
EffectiveFrequency	float	Observing frequency, in Hz
BeamSize	int	Beam size, in arcsec
NChannels	int	Total number of pixels in the instrument
Gain	float array	1D array with relative gains (flat field)
Offsets	float array	relative (X,Y) offsets, in arcsec
Channel_Sep	float array	matrix of channel to channel separations, in arcsec
TransmitionCurve	float array	
Flags	int array	Flag value for each channel (0 = unflagged)
RefChannel	int	Reference channel number
NUsedChannels	int	Number of channels in use for this observation
UsedChannels	int array	List of channels in use for this observation
CurrChanList	int array	Current list of channel numbers

Table 6.2: Attributes of a Telescope object

Name	Type	Description
Name	str	Telescope name, e.g. APEX-12m
Diameter	float	Antenna diameter, in m
Latitude	float	Latitude, in deg
Longitude	float	Longitude, in deg
Elevation	float	Elevation, in m

*Note:* for observations performed with wobbler switching, pairs of ON–OFF integrations are extracted from the Wobbler\_Sta attribute, and the phase differences are computed. By default, after reading, only the differentiated signals are stored in the Data attribute. The user can specify the phase number in the read command, in order to get only the 'ON' or the 'OFF' data.

### 6.2.2 DataAna

On top of the DataEntity, the DataAna layer defines additional attributes, related to statistics and flagging of the data. They are listed in Table 6.5.



Table 6.3: Attributes of the ScanParam object

Name	Type	Description
ScanNum	int	Scan number
ScanType	string	Scan type, e.g. 'FOCUS-Z
ScanMode	string	Scan mode, e.g. 'RASTER'
ScanDir	string	Scanning direction
Line_Len	float	Line length for a raster, in arcsec
Line_Ysp	float	Y-step between lines in a raster, in arcsec
Az_Vel	float	Scanning speed in Az, in arcsec/s
Object	string	Target name
Basis	tuple	Pair of strings describing basis frame - e.g. ('RA- - -SFL', 'DEC--SFL')
Coord	tuple	Target coordinates in basis frame
Date_Obs	string	Date of observation
Equinox	float	Equinox
Nula, Nule	floats	X, Y pointing settings at scan start
Colstart	float	Focus-Z setting at scan start
DeltaCA, DeltaIE	floats	Accumulated pointing corrections CA and IE
NObs	int	Number of subscans
SubscanNum	int list	Subscans numbers
SubscanIndex	int array	Integration numbers at subscans starts and ends
SubscanEpo	float array	Epochs of subscans starts, in year
SubscanTime	float array	LST times of subscans starts, in s
SubscanType	string list	Types of subscans - e.g. 'ON', or 'REF'
WobUsed	int	Boolean: is a wobbler used?
WobCycle	float	Wobbler period, in s
WobblerPos	float array	Wobbler positions, in arcsec
WobThrow	float	Wobbler throw, in arcsec
WobblerSta	string list	Wobbler status
Nodding_Sta	int array	Nodding status
WobMode	string	Wobbler mode, e.g. 'SQUARE'
AddLonWT	int	Wobbler throw to be added in Az, in arcsec
AddLatWT	int	Wobbler throw to be added in El, in arcsec
OnOffPairs	int list	List of pairs of integration numbers (if wobbler)
Nint	int	Number of integrations
Baslon, Baslat	float arrays	Absolute coordinates in basis frame, in deg
Track_Az, Track_El	float arrays	Tracking errors in Az and El, in arcsec
Lon, Lat	float arrays	Offsets w.r.t. the source in Az and El, in deg
FocX, FocY, FocZ	float arrays	Subreflector positions in X, Y, Z, in mm
PhiX, PhiY	float arrays	Subreflector rotation angles in X and Y, in deg
Az, El	float arrays	Absolute coordinates in Az, El, in deg
Lonpole, Latpole	float array	Coordinates in user frame of basis pole
Rot	float array	Rotation angle between user and basis frames, in deg
MJD	float array	Timestamps in MJD, in days

Table 6.4: Other attributes of a DataEntity object

Name	Type	Description
FileName	string	Input file name
RefGain	float	Frontend gain/attenuation factor
JyPerCount	float	Counts to Jy conversion factor
Data	float array	Current version of the data
DataBackup	float array	Previous version of the data
DataWeights	float array	Relative weights of the datapoints
DataFlags	array	Flagging of individual datapoints (0 = unflagged)
CorMatrix	float array	Channel to channel correlation matrix
FFCF_Gain	float array	1D array of relative gains (flat field) derived from skynoise
FFCF_CN	float array	Channel to channel correlated skynoise
SkyNoise	float array	Skynoise present in the signal

Table 6.5: Other attributes of a DataAna object

Name	Type	Description
ChanMean	float array	Mean values of signal per channel
ChanRms	float array	R.M.S of signal per channel
ChanMed	float array	Median values of signal per channel
ChanMean_s	float array	Mean values of signal per channel and per subscan
ChanRms_s	float array	R.M.S. of signal per channel and per subscan
ChanMed_s	float array	Median values of signal per channel and per subscan
flagValue	int	Current default flag value when calling a flagging routine
flagValueList	int list	Allowed values for flagging

### 6.2.3 Map

Finally, any kind of observation is stored in **BoA** in a Map object, that defines many methods for data reduction (see the Appendix for reference). It also contains an attribute called 'Map', of class Image, where the results of a map-making routine are stored.

### 6.2.4 Storing a data object

At any time during a **BoA** session, the user can dump the content of the current data object to a file. It can later be loaded again into **BoA**, in order to continue with the data reduction. This is done with:

```
boa> dump()
boa< I: current data successfully written to BoaData.sav
```

or:

```
boa> dump('myMap.data')
boa< I: current data successfully written to myMap.data
```

to give another filename than the default BoaData.sav. Then to reload the data object, one has to do:

```
boa> dd = newRestoreData()
```

**Note:** it is not possible in its present state to apply this restore method to the default *data* object. Therefore, after reloading a data object to a new variable (*dd* in the above example), one has to use the extended syntax (Chapter ...) instead of the abbreviations defined in BoaShortcuts.py.

## 6.3 Data output

### 6.3.1 Converting the raw data

**BoA** provides a procedure to convert an MB-FITS file to a FITS file with the same format as for MAMBO-ABBA data. The aim of this procedure is to be able to compare the results of a data reduction performed with **BoA** with those obtained with existing packages (e.g. NIC, MOPSI). *Note: This procedure has not been extensively tested recently...*

### 6.3.2 Saving a map

Once a mapping observation has been read in and processed with **BoA**, the user can store the results, i.e. a map in sky coordinates, in FITS file with standard 2D FITS images, including header with World Coordinate System (WCS) informations. This is done with the following command:

```
boa> data.writeFits()      # default file name: boaMap.fits
boa> data.writeFits('LABOCA_1234.fits') # give a file name
```

The resulting FITS file will contain three images, displaying the Intensity, the Rms, and the Weights of the current map. The content of each image is identified by the FITS keyword EXTNAME.

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## 7. DEVELOPMENT

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### 7.1 Basic programming rules

### 7.2 Adding classes

### 7.3 Adding methods

### 7.4 Adding Fortran90 code

*FB040510*

#### **General**

We are using Fortran 90/95 subroutines, wrapped to be called from python using the f2py package. This is because f90 code executes much faster than python scripts. There are some subtleties to pay attention to when wrapping fortran code, else you will add large overheads from the py-f90 interface, as arrays are copied and reindexed. For an introduction to F90/95 (only minor differences between the two), I recommend the compact and rather comprehensive (and free!) “Fortran 90 course notes”<sup>1</sup> by AC Marshall from the University of Liverpool. It contains all you probably need to know. I wrote a simple fortran method in BoA/fortran/BoaTest1.f90 to illustrate some basic features and give you a chance to test the wrapper without **BoA**. Look at its header for details. For an online F90/95 language reference<sup>2</sup> the best I found is at the NCSA resources page, describing IBM’s XL Fortran for AIX 8.1 – which is close to the Intel compiler.

#### **F90 in BoA**

For BoA our general idea is to have one f90.so extension module, which includes all the f90 methods (called subroutines and functions in fortran). This is necessitated by that the f90.data module, which contains much of a scans data, is connected (through an “use data”) to the other f90 program modules, and therefore they all need to be linked together.

The f90 methods may be split into different modules (classes) for convenience. We now have the first operational modules BoaF1.f90, BoaChannelAnalyser.f90, BoaBaseLine.f90, and the data module BoaData.f90. Each module may include any number of subroutines or functions. The data module BoaData.f90 is like a common block that contains all the data which does not change during data reduction. All data which does change is passed to the fortran subroutines as call arguments.

---

<sup>1</sup><http://math.nist.gov/WMitchell/f90course/CourseNotes.pdf>

<sup>2</sup>[http://www.ncsa.uiuc.edu/UserInfo/Resources/Hardware/IBMp690/IBM/usr/share/man/info/en\\_US/xlf/html/lr02.HTM#CONTENT](http://www.ncsa.uiuc.edu/UserInfo/Resources/Hardware/IBMp690/IBM/usr/share/man/info/en_US/xlf/html/lr02.HTM#CONTENT)

The `BoaData.f90` (`f90.data` from python) module is filled in `BoaDataEntity.FillF90`. It must be refilled if you change data object, else the fortran methods will work on a different scan. This re-filling must be implemented still. Currently the `f90.data` is only filled upon read of a new data file.

The CVS directory `BoA/fortran` contains the fortran source code. You will need to wrap/compile the **BoA** modules on your local system (see below), since it links to local libraries that have no standard address. This will create the extension module `f90.so` which you import to `BoA`. From the CVS directory `BoA` start **BoA**, then

```
>>> from fortran import f90
```

This is how to import any module from a subdirectory, which for this needs to include an empty file `__init__.py`

The python script `fortran/ftest.py` contains a series of calls to the fortran subroutines. To run it:

```
>>> read() # read in some scan
>>> op()   # open plot device
[enter]
>>> execfile('ftest.py') # start the script
```

which is followed with lots of output. To illustrate the use of new python methods that use fortran, you find `BoA/TestFB.py`, which you run like `ftest.py`. It goes through a number of data reduction steps and plots the data.

### Wrapping F90 code with f2py

To wrap the `f90` modules to produce `f90.so`:

```
ifc -c -w svd.f90
f2py -c -m f90 BoaData.f90 BoaF1.f90 BoaChannelAnalyser.f90 BoaBaseLine.f90 s
    or on some installations alternatively:
f2py -c --fcompiler=intel -m f90 BoaData.f90 BoaF1.f90 BoaChannelAnalyser.f90
```

The first command recompiles the `svd.o`. On the `f2py` line there are some diagnostic options you may add if you debug your code:

```
-DF2PY_REPORT_ATEXIT : gives time statistics upon exit from python.
-DF2PY_REPORT_ON_ARRAY_COPY=1000 : reports when the f2py interface copies an
-DNUMARRAY : must be used for numarray support. Default is Numeric.
```

If the wrapping fails, one of the following may be wrong:

1. You have not initiated the `ifc` compiler properly. In your shell initialization file (e.g. `.cshrc` for `tcsh`) you need

```
if (-e /opt/intel/compiler60/ia32/bin/ifcvars.sh) then
    source /opt/intel/compiler60/ia32/bin/ifcvars.csh
endif
```

or something equivalent.

2. Your python path does not include the intel fortran compiler:

```
setenv PYTHONPATH " ./opt/intel/compiler60/ia32/lib/:
                  /usr/local/lib/python2.3:
                  /usr/local/lib/python2.3/site-packages:
                  /home/bertoldi/bin:
                  /opt:
                  /usr/lib"
```

3. You use an old version of f2py.

```
<fortran> f2py -version
2.39.235_1644
```

Once you have successfully imported f90 in BoA, you can inquire about the use of a given method by typing

```
print f90.f1.NAME.__doc__
```

Fortran attributes are called f90.data.name\_of\_attribute. To inquire which ones are available:

```
boa> print f90.data.__doc__
el - 'f'-array(218)
track_el - 'f'-array(218)
ffcf_gain - 'f'-array(120)
subscan_time - 'f'-array(4)
az_p - 'f'-array(109,3)
lst - 'f'-array(218)
lon_p - 'f'-array(109,3)
track_az - 'f'-array(218)
lat - 'f'-array(218)
az - 'f'-array(218)
lat_p - 'f'-array(109,3)
lst_p - 'f'-array(109,3)
array_gain - 'f'-array(120)
lon - 'f'-array(218)
ffcf_cn - 'f'-array(120)
ut_p - 'f'-array(109,3)
nodding_sta - 'i'-array(218)
subscan_index - 'i'-array(4)
subscan_num - 'i'-array(4)
weights - 'f'-array(0), not allocated
el_p - 'f'-array(109,3)
ut - 'f'-array(218)
wobbler_pos - 'f'-array(218)
```

They are filled in in BoABusiness.py: BoAB.FillF90

### Use f90 methods in BoA

To call a fortran method, here an example:

```
compressed_array, nmax = f90.f1.compress(array, flag_array, 0)
```

Two objects are returned as a tuple, an array and an integer. They both are not in the call argument list, they are hidden to python, but are listed in the f90 code call argument list – have a look at the source code.

### Limitations

This particular example illustrates one of the limitations of wrapping f90 code: you cannot return an array with a length that is determined upon execution. The wrapper needs to specify the size of an array somehow. It does not have to be fixed, but specified through the size of an input attribute at least. In this example we try to return an array that is a compression of the input array, determined by the condition that the corresponding flag is 0. The trick to still do this here is to return a compressed\_array with the same size as array, plus an integer telling the size of the compressed array, so that the final answer is compressed\_array[0:nmax].

### Fortran vs. C-contiguous

If a Numeric array is proper-contiguous and has a proper type then it is directly passed to the wrapped Fortran function. Otherwise, an element-wise copy of an input array is made and the copy, being proper-contiguous and with proper type, is used as an array argument. There are two types of proper-contiguous Numeric arrays: Fortran-contiguous arrays when data is stored column-wise, i.e. indexing of data as stored in memory starts from the lowest dimension; C-contiguous when data is stored row-wise, i.e. indexing of data as stored in memory starts from the highest dimension. For one-dimensional arrays these notions coincide. To transform input arrays to column major storage order before passing them to Fortran routines, one may use the function `as_column_major_storage(<array>)` that is provided by all F2PY generated extension modules, such as the BoA f90. If you call a fortran method repeatedly with the same input array, you should convert the array first to avoid conversion by the wrapper interface on each call – which could dominate the execution time here. If you add the option `-DF2PY_REPORT_ON_ARRAY_COPY=1000` when wrapping, you will be informed on each copy that the wrapper interface performs. The option `-DF2PY_REPORT_ATEXIT` gives an execution time summary upon exit that splits up the time used in fortran and in the interface. If the interface time is large or comparable to the fortran execution time, your code is not efficient because it copies arrays too often. Look at examples in BoABaseLine.py, e.g.:

```
Data = f90.as_column_major_storage(self.Data.Data_Red_p)
Flag = f90.as_column_major_storage(self.Data.Data_Flag_p)
...
for i_ch in ch_range: # loop over channels and phases
    for i_ph in ph_range:
        Data = f90.baseline.addpoly(Data, Poly, Mean, Rms, i_ph, i_ch)
```

The input arrays are copied once into fortran-contiguous arrays before the loop, so in the loop there is no overhead from copying. Note also the general scheme of calling a fortran method here: Data is in- and output argument.



## 7.5 Interfacing

### 7.5.1 ScientificPython-2.4.5

ScientificPython is a collection of Python modules that are useful for scientific computing. Almost all modules make extensive use of Numerical Python (NumPy, Numeric), which must be installed prior to Scientific Python. Scientific consist of about one dozen modules, which contain methods written in Python that may come handy, but may be slow. The following lists a number of them.

stat() statistics() command calculates the statistics for all the channels in the range. Using plotmean() plotrms() we can plot mean and RMS values of each channels. The examples are as shows below:

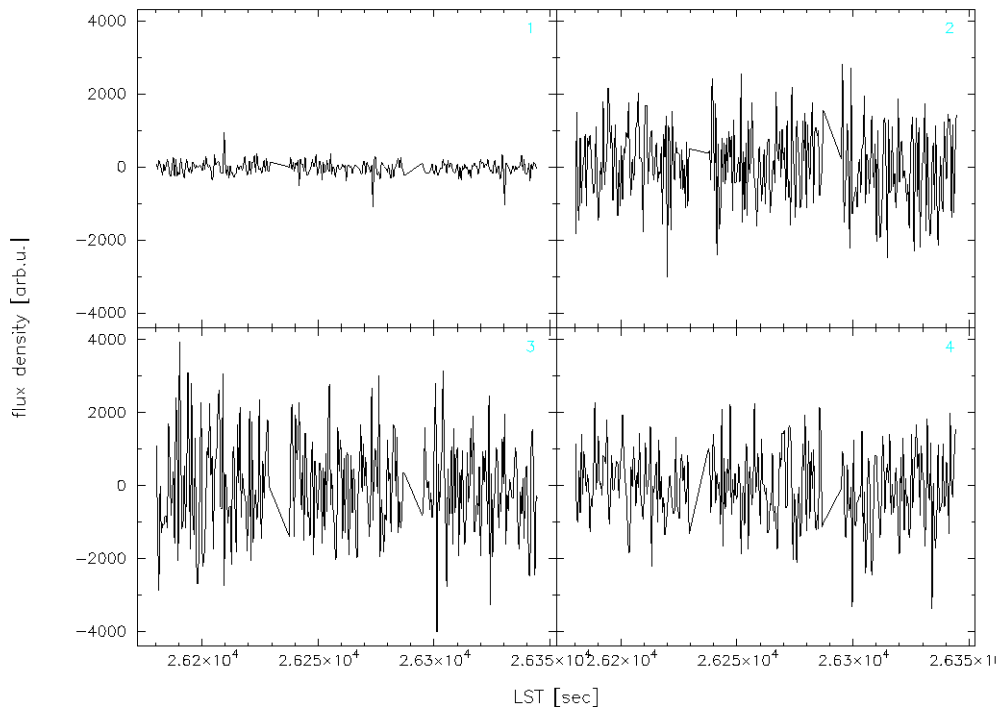


Figure 7.5.1: Plotting the Signal for channels in the range.

You need to import Numeric for Scientific. You can access the methods by importing the class or all methods:

```
>>> from Numeric import *
>>> import Scientific.Statistics
>>> Scientific.Statistics.median([1,2,3,5,6])
3.0
```

or alternatively

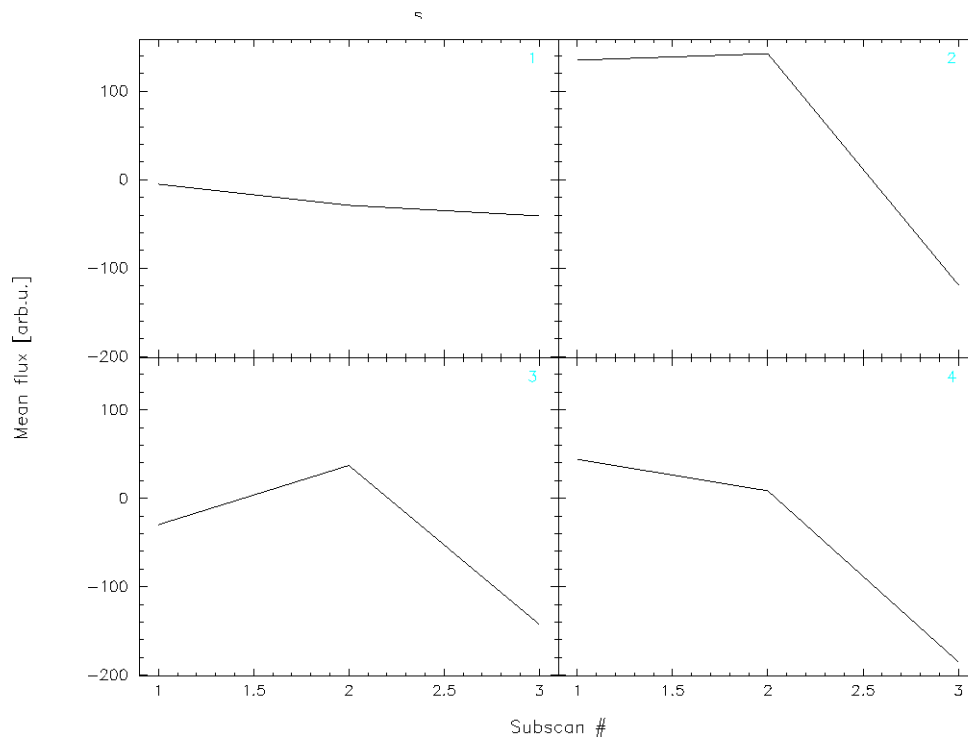


Figure 7.5.2: Plotting the Mean values of signal.

```
>>> from Scientific.Statistics import *
>>> median([1,2,3,5,6])
3.0
```

Available method in class Scientific.Statistics:

```
moment(data, order, about=None, theoretical=1)
mean(data)
weightedMean(data, sigma)
variance(data)
standardDeviation(data)
median(data)
mode(data)
normalizedMoment(data, order)
skewness(data)
kurtosis(data)
correlation(data1, data2)
```

There are also two classes for histograms:

```
Histogram
```

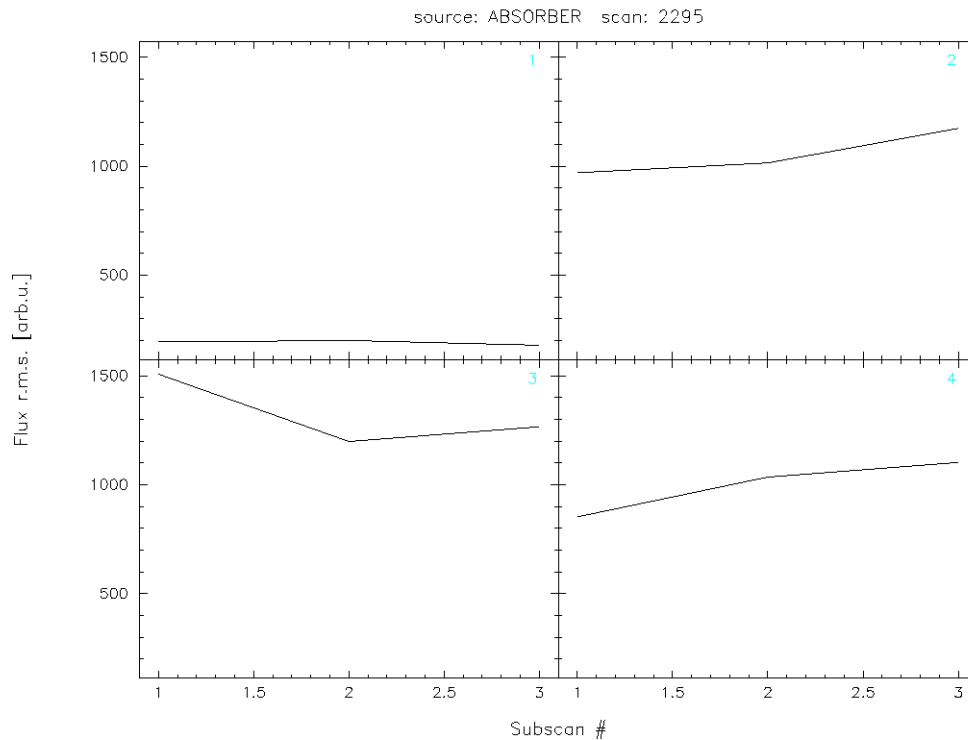


Figure 7.5.3: Plotting the RMS values of signal.

```
WeightedHistogram(Histogram)
```

The following explains only those Scientific methods which are useful for Boa. Consult the scripts or the (very sparse) documentation for more info.

### Scientific.Statistics.median

**Description:** Computes the median of a 1-d array.

**Example:**

```
>>> median([1, 2, 3, 5, 6])
3.0
```

### Scientific.Statistics.mean

**Description:** Returns the mean (average value) of a 1-d array.

**Example:**

```
>>> mean([1, 2, 3, 5, 6])
3.3999999999999999
```

### Scientific.Statistics.correlation

**Description:** Computes the correlation coefficient between two 1-dim arrays  $a$  and  $b$  according to

$$c_{ab} = \frac{\langle (a - \bar{a})(b - \bar{b}) \rangle}{\langle (a - \bar{a})^2 \rangle^{1/2} \langle (b - \bar{b})^2 \rangle^{1/2}} \quad (7.5.1)$$

### Example:

```
>>> correlation([1, 2, 3, 4, 5], [1, 2, 3, 4, 5])
1.0
>>> correlation([1, 2, 3, 4, 5], [1, 2, 3, 5, 5])
0.96476382123773219
>>> correlation([1, 2, 3, 4, 5], [5, 4, 3, 2, 1])
-1.0
```

### Scientific.Functions.LeastSquares

**Description:** General non-linear least-squares fit using the Levenberg-Marquardt algorithm and automatic derivatives. The parameter model specifies the function to be fitted. It will be called with two parameters: the first is a tuple containing all fit parameters, and the second is the first element of a data point (see below). The return value must be a number. Since automatic differentiation is used to obtain the derivatives with respect to the parameters, the function may only use the mathematical functions known to the module FirstDerivatives. The parameter parameter is a tuple of initial values for the fit parameters. The parameter data is a list of data points to which the model is to be fitted. Each data point is a tuple of length two or three. Its first element specifies the independent variables of the model. It is passed to the model function as its first parameter, but not used in any other way. The second element of each data point tuple is the number that the return value of the model function is supposed to match as well as possible. The third element (which defaults to 1.) is the statistical variance of the data point, i.e. the inverse of its statistical weight in the fitting procedure. The function returns a list containing the optimal parameter values and the chi-squared value describing the quality of the fit.

### Example:

```
>>> from Numeric import exp
>>> def f(param, t):
...     return param[0]*exp(-param[1]/t)
...
>>> data = [(100, 4.999e-8), (200, 5.307e+2),
            (300, 1.289e+6), (400, 6.559e+7)]
```

```
>>> print leastSquaresFit(f, (1e13,4700), data)
([8641551709749.7666, 4715.4677901570467], 1080.2526437958597)
```