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# **Numina Documentation**

*Release 0.21*

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**Jul 01, 2019**



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Welcome. This is the Documentation for Numina (version 0.21, date Jul 01, 2019),

Numina user guide: *Numina User Guide*

Numina pipeline creation guide: *Numina Pipeline Creation Guide*

Numina reference guide: *Numina Reference*.



This is Numina, the data reduction package used by the following GTC instruments: EMIR, MEGARA and FRIDA. Numina is distributed under GNU GPL, either version 3 of the License, or (at your option) any later version. See the file LICENSE.txt for details.

This guide is intended as an introductory overview of Numina and explains how to install and make use of the most important features. For detailed reference documentation of the functions and classes contained in the package, see the *Numina Reference*.

<p><b>Warning:</b> This “User Guide” is still a work in progress; some of the material is not organized, and several aspects of Numina are not yet covered sufficient detail.</p>
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## 1.1 Installation

### 1.1.1 Requirements

Python 2.7 or Python  $\geq 3.5$  is required. Additionally the following packages are required in order to work properly:

- setuptools
- six
- numpy
- scipy
- astropy
- PyYaml
- matplotlib
- python-dateutil

- `enum34` (Python 2.7 only)

Cython is required if you build the code from the development repository:

- `Cython`

The following packages are optional, for building documentation and testing:

- `sphinx` to build the documentation
- `pytest` for testing

### 1.1.2 Installing numina

The preferred methods of installation are `pip` or `conda`, using prebuilt packages.

#### Using pip

Run:

```
pip install numina
```

Pip will download all the required dependencies and a precompiled version of numina (if it exists for your platform) from PyPI.

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**Note:** If possible, pip will install a precompiled version of numina in wheel format. If such a version does not exist, pip will download and compile the source code. Numina has some portions of C and C++ code. You will need a C/C++ compiler such as `gcc` or `clang` (see *Building from source* below)

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Pip can install packages in different locations. You can use the `--user` option to install packages in your home directory.

Our recommended option is to perform isolated installations using `virtualenv` or `venv`. See *Numina Deployment with Virtualenv* for details.

**Warning:** Do not use `sudo pip` unless you *really really* know what you are doing.

#### Using conda

Numina packages for conda are provided in the `conda-forge` channel. To install the latest version of numina run:

```
conda update -c conda-forge numina
```

See *Numina Deployment with Conda* for details.

### 1.1.3 Building from source

You may end up building from source if there is not a precompiled version of numina for your platform or if you are doing development based on numina.



## Prerequisites

You will need a compiler suite and the development headers of Python and Numpy.

If you are building the development version of numina, you will also need Cython to translate Cython code into C/C++. If your source code is from a release, the translated files are included, and hence do not require Cython.

The released sources of numina can be downloaded from PyPI. If you require instead the development version, it can be checked out with:

```
git clone https://github.com/guaix-ucm/numina.git
```

## Building and installing

To build numina run:

```
python setup.py build
```

**Note:** In macOS Mojave, the compilation will fail unless the following environment variable is defined:

```
export MACOSX_DEPLOYMENT_TARGET=10.9
```

To install numina run:

```
python setup.py build
```

If you get an error about insufficient permissions to install, you are probably trying to access directories owned by root. Try instead:

```
python setup.py install --user
```

or perform the installation inside an isolated environment, such as conda or venv.

**Warning:** Do not `sudo python setup.py install` unless you really really know what you are doing.

### 1.1.4 Building the documentation

The Numina documentation is based on `sphinx`. With the package installed, the html documentation can be built from the `doc` directory:

```
$ cd doc
$ make html
```

The documentation will be copied to a directory under `build/sphinx`.

The documentation can be built in different formats. The complete list will appear if you type `make`

## 1.2 Numina Deployment with Virtualenv

`Virtualenv` is a tool to build isolated Python environments.

Since Python version 3.3, there is also a module in the standard library called *venv* with roughly the same functionality.

### 1.2.1 Create Virtual Environment

In order to create a virtual environment called e.g. *numinaenv* using *venv*:

```
python3 -m venv numinaenv
```

With *virtualenv* (in Python 2.7):

```
virtualenv numinaenv
```

### 1.2.2 Activate the Environment

Once the environment is created, you need to activate it. Just go to *bin/* folder created under *numinaenv* and source the script *activate*:

```
cd numinaenv/bin
source activate
(numinaenv) $
```

Notice that the prompt changes once you have activated the environment. To deactivate it, just type *deactivate*:

```
(numinaenv) $ deactivate
$
```

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**Note:** We are assuming that the user shell is *bash*. There are alternative *activate* scripts for *tcsh* and *fish* called *activate.csh* and *activate.fish*

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### 1.2.3 Numina Installation

After the environment activation, we can install *numina* with *pip*. This is the standard Python tool for package management. It will download the package and its dependencies, unpack everything and compile when needed:

```
(numinaenv) $ pip install numina
```

The requirements of *numina* will be downloaded and installed inside the virtual environment automatically.

You can also update *numina*, if your environment contains already a installed version:

```
(numinaenv) $ pip install -U numina
```

### 1.2.4 Test the installation

We can test the installation by running the *numina* command:

```
(numinaenv) $ numina
DEBUG: Numina simple recipe runner version 0.20
```

## 1.3 Numina Deployment with Conda

Conda was created with a target similar to [virtualenv](#), but extended its functionality to the management of packages in different languages.

You can install [miniconda](#) or [anaconda](#). The difference is that miniconda provides a light-weight environment and anaconda comes with lots of additional Python packages. By installing miniconda you reduce the amount of preinstalled packages in your system (after installing miniconda it is possible to install anaconda by executing `conda install anaconda`).

If you have updated the `$PATH` system variable during the miniconda or conda installation, you can call conda commands directly in the shell, like this:

```
bash$ conda info
```

If not, you will need to add the path to the command, like:

```
bash$ /path/to/conda/bin/conda info
```

In this guide we will write the commands without the full path, for simplicity.

Once conda is installed according to the corresponding miniconda or anaconda instructions, the steps to install numina under conda are:

### 1.3.1 Create a conda environment

With conda, environments are created in a centralised manner (under the subdirectory `./envs` in your conda tree):

```
conda create --name numinaenv
```

The Python interpreter used in this environment is the same version currently used by conda. You can select a different version with:

```
conda create --name numinaenv python=3.6
```

### 1.3.2 Activate the environment

With command:

```
conda activate numinaenv
```

which yields a different system prompt to the user:

```
(numinaenv) $
```

To exit the environment is enough to exit the terminal or run the following command:

```
(numinaenv) $ conda deactivate
```

### 1.3.3 Numina installation

After the environment activation, we can install numina using conda (we provide conda packages in the [conda-forge](#) channel):

```
(numinaenv) $ conda install -c conda-forge numina
```

We can also update numina, if your environment contains already a installed version:

```
(numinaenv) $ conda update numina
```

### 1.3.4 Test the installation

We can test the installation by running the numina command:

```
(numinaenv) $ numina  
DEBUG: Numina simple recipe runner version 0.20
```

## 1.4 Command Line Interface

The **numina** script is the interface with the pipelines It is called like this:

```
$ numina [global-options] comands [comand-options]
```

The **numina** script has several options:

- d, --debug**  
Debug enabled, increases verbosity.
- l filename**  
A file con configuration options for logging.

### 1.4.1 Options for run

The run subcommand processes the observing result with the appropriated reduction recipe.

It is called like this:

```
$ numina [global-options] run [comand-options] observation-result.yaml
```

- instrument 'name'**  
Name of one of the predefined instrument configurations.
- pipeline 'name'**  
Name of one of the predefined pipelines.
- requirements filename**  
File with the description of the parameters of the recipe.
- basedir path**  
File path used to resolve relative paths in the following options.
- datadir path**  
File path to the folder containing the pristine data to be processed.
- resultsdir path**  
File path to the directory where results are stored.
- workdir path**  
File path to the a directory where the recipe can write. Files in datadir are copied here.

**--cleanup**

Remove intermediate and temporal files created by the recipe.

**observing\_result** filename

Filename containing the description of the observation result.

## 1.4.2 Options for show-instruments

The show-instruments subcommand outputs information about the instruments with available pipelines.

It is called like this:

```
$ numina [global-options] show-instruments [options]
```

**-o, --observing-modes**

Show names and keys of Observing Modes in addition of instrument information.

**name**

Name of the instruments to show. If empty show all instruments.

## 1.4.3 Options for show-modes

The show-modes subcommand outputs information about the observing modes of the available instruments.

It is called like this:

```
$ numina [global-options] show-modes [options]
```

**-i, --instrument** name

Filter modes by instrument name.

**name**

Name of the observing mode to show. If empty show all observing modes.

## 1.4.4 Options for show-recipes

The show-recipes subcommand outputs information about the recipes of the available instruments.

It is called like this:

```
$ numina [global-options] show-recipes [options]
```

**-i, --instrument** name

Filter recipes by instrument name.

**-m, --mode**

Filter recipes by observing mode.

**name**

Name of the recipe to show. If empty show all recipes.



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## Numina Pipeline Creation Guide

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This guide is intended as an introductory overview of pipeline creation with Numina. For detailed reference documentation of the functions and classes contained in the package, see the *Numina Reference*.

**Warning:** This “Pipeline Creation Guide” is still a work in progress; some of the material is not organized, and several aspects of Numina are not covered in sufficient detail yet.

## 2.1 Numina Pipeline Concepts

### 2.1.1 Instrument

### 2.1.2 Observing Modes

Each Instrument has a list of predefined types of observations that can be carried out with it. Each Observing Mode is defined by:

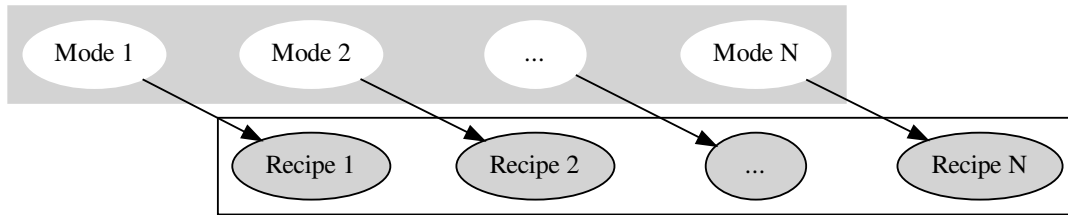
- The configuration of the Telescope
- The configuration of the Instrument
- The type of processing required by the images obtained during the observation

Some of the observing modes of a Instrument are **Scientific**, that is, modes devoted to obtain data to perform scientific analysis. Other modes are devoted to **Calibration**; these modes produce data required to correct the scientific images from the effects of the Instrument, the Telescope and the atmosphere.

### 2.1.3 Recipes

A recipe is a method to process the images obtained in a particular observing mode. Recipes in general require (as inputs) the list of raw images obtained during the observation. Recipes can require other inputs (calibrations), and those inputs can be the outputs of other recipes.

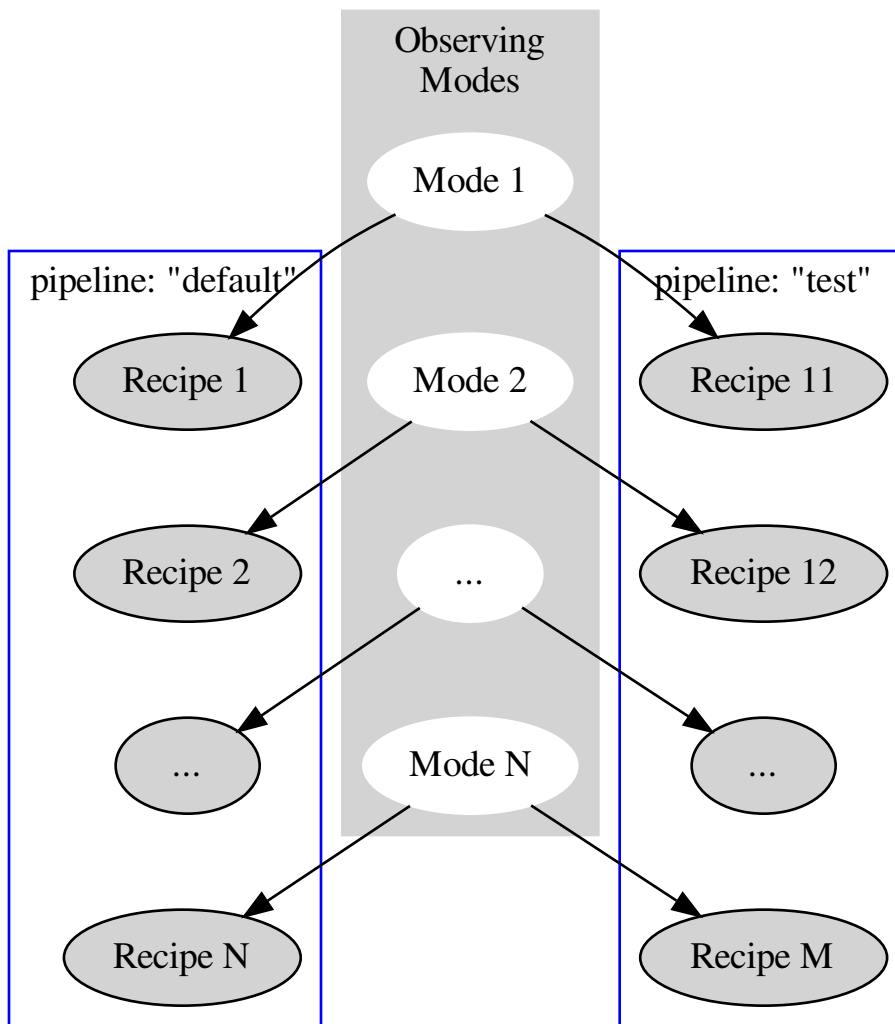
Images obtained in a particular mode are processed by one recipe.



### 2.1.4 Pipelines

A pipeline represents a particular mapping between the observing modes and the reduction algorithms that process each mode. Each instrument has at least one pipeline called *default*. It may have other pipelines for specific purposes.





### 2.1.5 Products, Requirements and Data Types

A recipe announces its required inputs as *Requirement* and its outputs as *Result*.

Both Results and Requirements have a name and a type. Types can be plain Python types or defined by the developer.

### 2.1.6 Format of the input files

The default format of the input and output files is [YAML](#), a data serialization language.

## Format of the Observation Result file

This file contains the result of an observation. It represents an *ObservationResult* object.

The contents of the object are serialized as a dictionary with the following keys:

- id: not required, integer, defaults to 1** Unique identifier of the observing block
- instrument: required, string** Name of the instrument, as it appears in the instrument file (see below)
- mode: required, string** Name of the observing mode
- children: not required, list of integers, defaults to empty list** Identifications of nested observing blocks
- frames: required, list of file names** List of raw images

```
id: 21
instrument: EMIR
mode: nb_image
children: []
frames:
- r0121.fits
- r0122.fits
- r0123.fits
- r0124.fits
- r0125.fits
- r0126.fits
- r0127.fits
- r0128.fits
- r0129.fits
- r0130.fits
- r0131.fits
- r0132.fits
```

## Format of the requirement file (version 1)

```
version: 1
products:
  EMIR:
    - {id: 1, content: 'file1.fits', type: 'MasterFlat', tags: {'filter': 'J'}, ob: 200}
    - {id: 4, content: 'file4.fits', type: 'MasterBias', tags: {'readmode': 'cds'}, ob: 400}
  MEGARA:
    - {id: 1, content: 'file1.fits', type: 'MasterFlat', tags: {'vph': 'LR1'}, ob: 1200}
    - {id: 2, content: 'file2.yml', type: 'TraceMap', tags: {'vph': 'LR2', 'readmode': 'fast'}, ob: 1203}
requirements:
  EMIR:
    default:
      TEST6:
        pinhole_nominal_positions: [ [0, 1], [0, 1] ]
        box_half_size: 5
      TEST9:
        median_filter_size: 5
  MEGARA:
    default:
      mos_image: {}
```

## Format of the requirement file

**Warning:** This section documents a deprecated format

Deprecated since version 0.14.0.

This file contains configuration parameters for the recipes that are not related to the particular instrument used.

The contents of the file are serialized as a dictionary with the following keys:

**requirements: required, dictionary** A dictionary of parameter names and values.

**logger: optional, dictionary** A dictionary used to configure the custom file logger

```
requirements:
  master_bias: master_bias-1.fits
  master_bpm: bpm.fits
  master_dark: master_dark-1.fits
  master_intensity_ff: master_flat.fits
  nonlinearity: [1.0, 0.0]
  subpixelization: 4
  window:
    - [800, 1500]
    - [800, 1500]
logger:
  logfile: processing.log
  format: "%(asctime)s - %(name)s - %(levelname)s - %(message)s"
  enabled: true
```

## 2.2 Numina Pipeline Example

This guide is intended as an introductory overview of the creation of instrument reduction pipelines with Numina. For detailed reference documentation of the functions and classes contained in the package, see the [Numina Reference](#).

**Warning:** This “Pipeline Creation Guide” is still a work in progress; some of the material is not organized, and several aspects of Numina are not yet covered sufficient detail.

### 2.2.1 Execution environment of the Recipes

Recipes have different execution environments. Some recipes are designed to process observing modes required for the observation. These modes are related to visualization, acquisition and focusing. The Recipes are integrated in the GTC environment. We call these recipes the **Data Factory Pipeline**, (*DFP*).

Other group of recipes are devoted to scientific observing modes: imaging, spectroscopy and auxiliary calibrations. These Recipes constitute the **Data Reduction Pipeline**, (*DRP*). The software is meant to be standalone, users shall download the software and run it in their own computers, with reduction parameters and calibrations provided by the instrument team.

Users of the DRP will use the simple Numina CLI. Users of the DFP shall interact with the software through the GTC Inspector.

## 2.2.2 Instrument Reduction Pipeline Example

In the following sections we create an Instrument Reduction Pipeline for an instrument name *CLODIA*.

In order to make a new Instrument Reduction Pipeline visible to Numina and the GTC Control System you have to create a full Python package that will contain the reduction recipes, data types and other processing code.

The creation of Python packages is described in detail (for example) in the [Python Packaging User Guide](#).

Then, we create a Python package called *clodiadrp* with the following structure (we ignore files such as README or LICENSE as they are not relevant here):

```
clodiadrp
|-- clodiadrp
|   |-- __init__.py
|-- setup.py
```

From here the steps are:

1. Create a configuration yaml file.
2. Create a loader file.
3. Link the *entry\_point* option in the *setup.py* with the loader file.
4. Create the Pipeline's Recipes.

In the following we will continue with the same example as previously.

### Configuration File

The configuration file contains basic information such as:

- the list of modes of the instrument
- the list of recipes of the instrument
- the mapping between recipes and modes.

In this example, we assume that CLODIA has three modes: **Bias**, **Flat** and **Image**. The first two modes are used for pedestal and flat-field illumination correction. The third is the main scientific mode of the instrument.

Create a new yaml file in the root folder named *drp.yaml*.

```
name: CLODIA
configurations:
  default: {}
modes:
  -key: bias
    name: Bias
    summary: Bias mode
    description: >
      Full description of the Bias mode
  -key: flat
    name: Flat
    summary: Flat mode
    description: >
      Full description of the Flat mode
  -key: image
    name: Image
    summary: Image mode
```

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```

description: >
  Full description of the Image mode
pipelines:
  default:
    version: 1
    recipes:
      bias: clodiadrp.recipes.recipe
      flat: clodiadrp.recipes.recipe
      image: clodiadrp.recipes.recipe

```

The entry *modes* contains a list of the observing modes of the instrument. There are three: Bias, Flat and Image. Each entry contains information about the mode. A *name*, a short *summary* and a multi-line *description*. The field *key* is used to map the observing modes and the *recipes*, so *key* has to be unique and equal to only one value in each *recipes* block under *pipelines*.

The entry *pipelines* contains only one pipeline, called *default* by convention. The *pipeline* contains recipes, each related to one observing mode by means of the filed *key*. For the moment we haven't developed any recipe, so the value of each key (*clodiadrp.recipes.recipe*) doesn't exist yet.

---

**Note:** This file has to be included in *package\_data* inside *setup.py* to be distributed with the package, see [Installing Package Data](#) for details.

---

## Loader File

Create a new loader file in the root folder named *loader.py* with the following information:

```

import numina.core

def drp_load():
    """Entry point to load CLODIA DRP."""
    return numina.core.drp_load('clodiadrp', 'drp.yaml')

```

## Create entry point

Once we have created the *loader.py* file, the only thing we have to do is to make CLODIA visible to Numina/GCS. To do so, just modify the *setup.py* file to add an entry point.

```

from setuptools import setup

setup(name='clodiadrp',
      entry_points = {
          'numina.pipeline.1': ['CLODIA = clodiadrp.loader:drp_load'],
      },
)

```

Both the Numina CLI tool and GCS check this particular entry point. They call the function provided by the entry point. The function *drp\_load()* reads and parses the YAML file and creates an object of class *InstrumentDRP* for each recipes it finds. These objects are used by Numina CLI and GCS to discover the available Instrument Reduction Pipelines.

At this stage, the file layout is as follows:

```
clodiadrp
|-- clodiadrp
|   |-- __init__.py
|   |-- loader.py
|   |-- drp.yaml
|-- setup.py
```

**Note:** In fact, it is not necessary to use a YAML file to contain the Instrument information. The only strict requirement is that the function in the entry point ‘numina.pipeline.1’ must return a valid *InstrumentDRP* object. The use of a YAML file and the *drp\_load()* function is only a matter of convenience.

---

## Recipes Creation

We haven’t created any reduction recipe yet. As a matter of organization, we suggest to create a dedicated subpackage for recipes *clodiadrp.recipes* and a module for each recipe. The file layout is:

```
clodiadrp
|-- clodiadrp
|   |-- __init__.py
|   |-- loader.py
|   |-- drp.yaml
|   |-- recipes
|   |   |-- __init__.py
|   |   |-- bias.py
|   |   |-- flat.py
|   |   |-- image.py
|-- setup.py
```

Recipes must provide three things: 1) a description of the inputs of the recipe; 2) a description of the products of the recipe and 3) a *run* method which is in charge of executing the processing. Additionally, all Recipes must inherit from *BaseRecipe*.

We start with a simple *Bias* recipe. Its purpose is to process images previously taken in *Bias* mode, that is, a series of pedestal images. The recipe will receive the result of the observation and return a master bias image.

```
from numina.core import Result, Requirement
from numina.core import DataFrameType
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe

class Bias(BaseRecipe):                                     (1)

    obresult = Requirement(ObservationResultType, "Observation Result") (2)
    master_bias = Result(DataFrameType)                    (3)

    def run(self, rinput):                                  (4)

        # Here the raw images are processed
        # and a final image myframe is created

        result = self.create_result(master_bias=myframe) (5)
        return result
```

1. Each recipe must be a class derived from *BaseRecipe*

2. This recipe only requires the result of the observation. Each requirement is an object of the `Requirement` class or any subclass of it. The type of the requirement is `ObservationResultType`, representing the result of the observation.
3. This recipe only produces one result. Each product is an object of `Result` class. The type of the product is given by `DataFrameType`, representing an image.
4. Each recipe must provide a `run` method. The method has only one argument that collects the values of all inputs declared by the recipe. In this case, `rinput` has a member named `obresult` and can be accessed through `rinput.obresult` which belongs to `ObservationResult` class.
5. The recipe must return an object that collects all the declared products of the recipe, of `RecipeResult` class. This is accomplished internally by the `create_result` method. It will raise a run time exception if any of the declared products are not provided.

We can now create the `Flat` recipe (inside `flat.py`). This recipe has two requirements, the observation result and a master bias image (flat-field images require bias subtraction).

```

from numina.core import Result, Requirement
from numina.core import DataFrameType
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe

class Flat(BaseRecipe):

    obresult = Requirement(ObservationResultType, "Observation Result") (1)
    master_bias = Requirement(DataFrameType, "Master Bias") (2)
    master_flat = Result(DataFrameType)

    def run(self, rinput): (3)

        # Here the raw images are processed
        # and a final image myframe is created

        result = self.create_result(master_flat=myframe) (4)
        return result

```

1. This recipe only requires the result of the observation. Each requirement is an object of the `Requirement` class or any subclass of it. The type of the requirement is `ObservationResultType`, representing the result of the observation.
2. It also requires a master bias image which belongs to `DataFrameType` class (represents an image).
3. In this case, `rinput` has two members: 1) `rinput.obresult` of `ObservationResult` class and 2) a `rinput.master_bias` of `DataFrame` class
4. The arguments of `create_result` must be the same names used in the product definition.

Finally, the recipe for `Image` mode reduction (inside `image.py`) has three requirements, the observation result, a master bias and a master flat images

```

from numina.core import Result, Requirement
from numina.core import DataFrameType
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe

class Image(BaseRecipe):

    obresult = Requirement(ObservationResultType)
    master_bias = Requirement(DataFrameType)

```

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```

master_flat = Requirement(DataFrameType)
final = Result(DataFrameType)

def run(self, rinput):
    # Here the raw images are processed
    # and a final image myframe is created

    result = self.create_result(final=myframe)
    return result

```

1. In this case, *rinput* will have three members *rinput.obresult* of *ObservationResult* class, *rinput.master\_bias* of *DataFrame* class and *rinput.master\_flat* of *DataFrame* class.

**Note:** It is not strictly required that the requirements and products names are consistent between recipes, although it is highly recommended.

Now we must update *drp.yaml* to insert the full name of the recipes (package and class), as follows

```

name: CLODIA
configurations:
  default: {}
modes:
  -key: bias
    name: Bias
    summary: Bias mode
    description: >
      Full description of the Bias mode
  -key: flat
    name: Flat
    summary: Flat mode
    description: >
      Full description of the Flat mode
  -key: image
    name: Image
    summary: Image mode
    description: >
      Full description of the Image mode
pipelines:
  default:
    version: 1
    recipes:
      bias: clodiadrp.recipes.bias.Bias
      flat: clodiadrp.recipes.flat.Flat
      image: clodiadrp.recipes.image.Image

```

## Specialized data products

There is some information that is missing of our current setup. The products of some recipes are the inputs of others. The master bias created by *Bias* is the input that *Flat* and *Image* require. To represent this situation we use specialized data products. We start by adding a new module *products*:



```

clodiadrp
|-- clodiadrp
|   |-- __init__.py
|   |-- loader.py
|   |-- products.py
|   |-- drp.yaml
|   |-- recipes
|   |   |-- __init__.py
|   |   |-- bias.py
|   |   |-- flat.py
|   |   |-- image.py
|-- setup.py

```

We have two types of images that are products of recipes that can be required by other recipes: **master bias** and **master flat**. We represent this by creating two new types derived from `DataFrameType` (because the new types are images) and `DataProductMixin` (because the new types are products that must be handled by both Numina CLI and GTC Control system) classes.

```

from numina.types.frame import DataFrameType
from numina.types.product import DataProductMixin

class MasterBias(DataFrameType, DataProductMixin):
    pass

class MasterFlat(DataFrameType, DataProductMixin):
    pass

```

Now we must modify our recipes as follows. First *Bias*

```

from numina.core import Result, Requirement
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe
from clodiadrp.products import MasterBias    (1)

class Bias(BaseRecipe):

    obresult = Requirement(ObservationResultType)
    master_bias = Result(MasterBias)          (2)

    ...                                     (3)

```

1. Import the new type *MasterBias*.
2. Declare that our recipe produces *MasterBias* images.
3. *run* method remains unchanged.

Then *Flat*:

```

from numina.core import Result, Requirement
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe
from clodiadrp.products import MasterBias, MasterFlat

class Flat(BaseRecipe):

    obresult = Requirement(ObservationResultType)

```

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```

master_bias = Requirement(MasterBias)          (1)
master_flat = Result(MasterFlat)              (2)
...                                           (3)

```

1. *MasterBias* is used as a requirement. This guaranties that the images provided here are those created by *Bias* and no other.
2. Declare that our recipe produces *MasterFlat* images.
3. *run* method remains unchanged.

And finally *Image*:

```

from numina.core import Result, Requirement
from numina.core import DataFrameType
from numina.types.obsresult import ObservationResultType
from numina.core.recipes import BaseRecipe
from clodiadrp.products import MasterBias, MasterFlat

class Image(BaseRecipe):

    obresult = Requirement(ObservationResultType)
    master_bias = Requirement(MasterBias)          (1)
    master_flat = Requirement(MasterFlat)          (2)
    final = Result(DataFrameType)                 (3)
    ...                                           (4)

```

1. *MasterBias* is used as a requirement. This guaranties that the images provided here are those created by *Bias* and no other.
2. *MasterFlat* is used as a requirement. This guaranties that the images provided here are those created by *Flat* and no other.
3. Declare that our recipe produces *Image* images.
4. *run* method remains unchanged.

## 2.3 DRP Data Types

Custom data types can be used as Requirements and Products by Recipes. New data types can be derived as follows.

### 2.3.1 Create a new DataType

New Data Types must derive from *numina.core.DataType* or one of its subclasses. In the constructor, we must declare the base type of the objects of this Data Product.

For example, a *MasterBias* Data Product is an image, so its base type is a *DataFrame*. A table of 2D coordinates will have a *numpy.ndarray* base type.

In general, we are interested in defining new DataTypes for objects that will contain information that will be used as inputs in different recipes. In this case, we must derive from *numina.core.DataProductType*.

As an example, we create a DataType that will store information about the trace of a spectrum. The information will be stored in Python *dict*.

```
class TraceMap(DataProductType):
    def __init__(self, default=None):
        super(TraceMap, self).__init__(dict, default)
```

### 2.3.2 Construction of objects

The input of a recipe is created by inspecting the Recipe Requirements. The Recipe Loader is in charge of finding an appropriated value for each requirement. The value is passed to *Requirement.convert*, that in turn calls *DataType.convert*. The default implementation just returns in input object unchanged.

### 2.3.3 Loading and Storage with the command line Recipe Loader

Each Recipe Loader can implement its own mechanism to store and load Data Products. The Command Line Recipe Loader uses text files in YAML format.

To define how a particular DataProduct is stored under the default Recipe Loader, two functions must be defined, a store function and a load function. Then these two functions must be registered with the global functions *numina.store.dump* and *numina.store.load*.

```
from numina.store import dump, load

from .products import TraceMap

@dump.register(TraceMap)
def dump_tracemap(tag, obj, where):

    filename = where.destination + '.yaml'

    with open(filename, 'w') as fd:
        yaml.dump(obj, fd)

    return filename

@load.register(TraceMap)
def load_tracemap(tag, obj):

    with open(obj, 'r') as fd:
        traces = yaml.load(fd)

    return traces
```

In this example, *tag* is an argument of type *TraceMap* and *obj* is of type *dict*.



**Release** 0.21

**Date** Jul 01, 2019

**Warning:** This “Reference” is still a work in progress; some of the material is not organized, and several aspects of Numina are not yet covered sufficient detail.

## 3.1 Numina modules

### 3.1.1 `numina.array` — Array manipulation

`numina.array.fixpix` (*data*, *mask*, *kind='linear'*)  
Interpolate 2D array data in rows

`numina.array.fixpix2` (*data*, *mask*, *iterations=3*, *out=None*)  
Substitute pixels in mask by a bilinear least square fitting.

`numina.array.numberarray` (*x*, *shape*)  
Return *x* if it is an array or create an array and fill it with *x*.

`numina.array.rebin` (*a*, *newshape*)  
Rebin an array to a new shape.

`numina.array.rebin_scale` (*a*, *scale=1*)  
Scale an array to a new shape.

`numina.array.subarray_match` (*shape*, *ref*, *sshape*, *sref=None*)  
Compute the slice representation of intersection of two arrays.

Given the shapes of two arrays and a reference point *ref*, compute the intersection of the two arrays. It returns a tuple of slices, that can be passed to the two images as indexes

**Parameters**

- **shape** – the shape of the reference array
- **ref** – coordinates of the reference point in the first array system
- **sshape** – the shape of the second array

**Param** sref: coordinates of the reference point in the second array system, the origin by default

**Returns** two matching slices, corresponding to both arrays or a tuple of Nones if they don't match

**Return type** a tuple

**Example**

```
>>> import numpy
>>> im = numpy.zeros((1000, 1000))
>>> sim = numpy.ones((40, 40))
>>> i, j = subarray_match(im.shape, [20, 23], sim.shape)
>>> im[i] = 2 * sim[j]
```

`numina.array.process_ramp` (*inp*[, *out=None*, *axis=2*, *ron=0.0*, *gain=1.0*, *nsig=4.0*, *dt=1.0*, *saturation=65631*])

New in version 0.8.2.

Compute the result 2d array computing slopes in a 3d array or ramp.

**Parameters**

- **inp** – input array
- **out** – output array
- **axis** – unused
- **ron** – readout noise of the detector
- **gain** – gain of the detector
- **nsig** – rejection level to detect glitched and cosmic rays
- **dt** – time interval between exposures
- **saturation** – saturation level

**Returns** a 2d array

**numina.array.background — Background estimation**

Background estimation

Background estimation following Costa 1992, Bertin & Arnouts 1996

`numina.array.background.background_estimator` (*bdata*)

Estimate the background in a 2D array

`numina.array.background.create_background_map` (*data*, *bsx*, *bsy*)

Create a background map with a given mesh size

**numina.array.blocks — Generation of blocks**

`numina.array.blocks.blk_1d` (*blk*, *shape*)

Iterate through the slices that recover a line.

This function is used by `blk_nd()` as a base 1d case.

The last slice is returned even if is lesser than `blk`.

**Parameters**

- **blk** – the size of the block
- **shape** – the size of the array

**Returns** a generator that yields the slices

`numina.array.blocks.blk_1d_short` (*blk*, *shape*)

Iterate through the slices that recover a line.

This function is used by `blk_nd_short()` as a base 1d case.

The function stops yielding slices when the size of the remaining slice is lesser than `blk`.

**Parameters**

- **blk** – the size of the block
- **shape** – the size of the array

**Returns** a generator that yields the slices

`numina.array.blocks.blk_coverage_1d` (*blk*, *size*)

Return the part of a 1d array covered by a block.

**Parameters**

- **blk** – size of the 1d block
- **size** – size of the 1d a image

**Returns** a tuple of size covered and remaining size

**Example**

```
>>> blk_coverage_1d(7, 100)
(98, 2)
```

`numina.array.blocks.blk_nd` (*blk*, *shape*)

Iterate through the blocks that cover an array.

This function first iterates trough the blocks that recover the part of the array given by `max_blk_coverage` and then iterates with smaller blocks for the rest of the array.

**Parameters**

- **blk** – the N-dimensional shape of the block
- **shape** – the N-dimensional shape of the array

**Returns** a generator that yields the blocks

### Example

```
>>> result = list(blk_nd(blk=(5,3), shape=(11, 11)))
>>> result[0]
(slice(0, 5, None), slice(0, 3, None))
>>> result[1]
(slice(0, 5, None), slice(3, 6, None))
>>> result[-1]
(slice(10, 11, None), slice(9, 11, None))
```

The generator yields blocks of size `blk` until it covers the part of the array given by `max_blk_coverage()` and then yields smaller blocks until it covers the full array.

#### See also:

`blk_nd_short()` Yields blocks of fixed size

`numina.array.blocks.blk_nd_short` (*blk*, *shape*)

Iterate through the blocks that strictly cover an array.

Iterate through the blocks that recover the part of the array given by `max_blk_coverage`.

#### Parameters

- **blk** – the N-dimensional shape of the block
- **shape** – the N-dimensional shape of the array

**Returns** a generator that yields the blocks

### Example

```
>>> result = list(blk_nd_short(blk=(5,3), shape=(11, 11)))
>>> result[0]
(slice(0, 5, None), slice(0, 3, None))
>>> result[1]
(slice(0, 5, None), slice(3, 6, None))
>>> result[-1]
(slice(5, 10, None), slice(6, 9, None))
```

In this case, the output of `max_blk_coverage` is (10, 9), so only this part of the array is covered

#### See also:

`blk_nd()` Yields blocks of `blk` size until the remaining part is smaller than `blk` and then yields smaller blocks.

`numina.array.blocks.block_view` (*arr*, *block*=(3, 3))

Provide a 2D block view to 2D array.

No error checking made. Therefore meaningful (as implemented) only for blocks strictly compatible with the shape of A.

`numina.array.blocks.blockgen` (*blocks*, *shape*)

Generate a list of slice tuples to be used by `combine`.

The tuples represent regions in an N-dimensional image.

#### Parameters

- **blocks** – a tuple of block sizes



- **shape** – the shape of the n-dimensional array

**Returns** an iterator to the list of tuples of slices

### Example

```
>>> blocks = (500, 512)
>>> shape = (1040, 1024)
>>> for i in blockgen(blocks, shape):
...     print i
(slice(0, 260, None), slice(0, 512, None))
(slice(0, 260, None), slice(512, 1024, None))
(slice(260, 520, None), slice(0, 512, None))
(slice(260, 520, None), slice(512, 1024, None))
(slice(520, 780, None), slice(0, 512, None))
(slice(520, 780, None), slice(512, 1024, None))
(slice(780, 1040, None), slice(0, 512, None))
(slice(780, 1040, None), slice(512, 1024, None))
```

`numina.array.blocks.blockgen1d` (*block*, *size*)

Compute 1d block intervals to be used by combine.

`blockgen1d` computes the slices by recursively halving the initial interval (0, size) by 2 until its size is lesser or equal than `block`

#### Parameters

- **block** – an integer maximum block size
- **size** – original size of the interval, it corresponds to a 0:size slice

**Returns** a list of slices

### Example

```
>>> blockgen1d(512, 1024)
[slice(0, 512, None), slice(512, 1024, None)]
```

`numina.array.blocks.max_blk_coverage` (*blk*, *shape*)

Return the maximum shape of an array covered by a block.

#### Parameters

- **blk** – the N-dimensional shape of the block
- **shape** – the N-dimensional shape of the array

**Returns** the shape of the covered region

### Example

```
>>> max_blk_coverage(blk=(7, 6), shape=(100, 43))
(98, 42)
```

### `numina.array.bpm` — Bad Pixel Mask interpolation

Fix points in an image given by a bad pixel mask

### `numina.array.combine` — Array combination

Different methods for combining lists of arrays.

`numina.array.combine.flatcombine` (*arrays, masks=None, dtype=None, scales=None, low=3.0, high=3.0, blank=1.0*)

Combine flat arrays.

#### Parameters

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays
- **blank** – non-positive values are substituted by this on output

**Returns** mean, variance of the mean and number of points stored

`numina.array.combine.generic_combine` (*method, arrays, masks=None, dtype=None, out=None, zeros=None, scales=None, weights=None*)

Stack arrays using different methods.

#### Parameters

- **method** (*PyCObject*) – the combination method
- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **zeros** –
- **scales** –
- **weights** –

**Returns** median, variance of the median and number of points stored

`numina.array.combine.mean` (*arrays, masks=None, dtype=None, out=None, zeros=None, scales=None, weights=None*)

Combine arrays using the mean, with masks and offsets.

Arrays and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the mean, `out[1]` the variance and `out[2]` the number of points used.

#### Parameters

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays

**Returns** mean, variance of the mean and number of points stored

### Example

```
>>> import numpy
>>> image = numpy.array([[1., 3.], [1., -1.4]])
>>> inputs = [image, image + 1]
>>> mean(inputs)
array([[ 1.5,  3.5],
       [ 1.5, -0.9]],
<BLANKLINE>
       [[ 0.5,  0.5],
       [ 0.5,  0.5]],
<BLANKLINE>
       [[ 2. ,  2. ],
       [ 2. ,  2. ]])
```

`numina.array.combine.median`(*arrays*, *masks=None*, *dtype=None*, *out=None*, *zeros=None*, *scales=None*, *weights=None*)

Combine arrays using the median, with masks.

Arrays and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the mean, `out[1]` the variance and `out[2]` the number of points used.

#### Parameters

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays

**Returns** median, variance of the median and number of points stored

`numina.array.combine.minmax`(*arrays*, *masks=None*, *dtype=None*, *out=None*, *zeros=None*, *scales=None*, *weights=None*, *nmin=1*, *nmax=1*)

Combine arrays using mix max rejection, with masks.

Inputs and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the mean, `out[1]` the variance and `out[2]` the number of points used.

#### Parameters

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays
- **nmin** –
- **nmax** –

**Returns** mean, variance of the mean and number of points stored

`numina.array.combine.quantileclip` (*arrays, masks=None, dtype=None, out=None, zeros=None, scales=None, weights=None, fclip=0.1*)

Combine arrays using the sigma-clipping, with masks.

Inputs and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the mean, `out[1]` the variance and `out[2]` the number of points used.

**Parameters**

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays
- **fclip** – fraction of points removed on both ends. Maximum is 0.4 (80% of points rejected)

**Returns** mean, variance of the mean and number of points stored

`numina.array.combine.sigmaclip` (*arrays, masks=None, dtype=None, out=None, zeros=None, scales=None, weights=None, low=3.0, high=3.0*)

Combine arrays using the sigma-clipping, with masks.

Inputs and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the mean, `out[1]` the variance and `out[2]` the number of points used.

**Parameters**

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **out** – optional output, with one more axis than the input arrays
- **low** –
- **high** –

**Returns** mean, variance of the mean and number of points stored

`numina.array.combine.sum` (*arrays, masks=None, dtype=None, out=None, zeros=None, scales=None*)

Combine arrays by addition, with masks and offsets.

Arrays and masks are a list of array objects. All input arrays have the same shape. If present, the masks have the same shape also.

The function returns an array with one more dimension than the inputs and with size (3, shape). `out[0]` contains the sum, `out[1]` the variance and `out[2]` the number of points used.

**Parameters**

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output

- **out** – optional output, with one more axis than the input arrays

**Returns** sum, variance of the sum and number of points stored

### Example

```
>>> import numpy
>>> image = numpy.array([[1., 3.], [1., -1.4]])
>>> inputs = [image, image + 1]
>>> sum(inputs)
array([[ [ 1.5,  3.5],
         [ 1.5, -0.9]],
       <BLANKLINE>
         [[ 0.5,  0.5],
          [ 0.5,  0.5]],
       <BLANKLINE>
         [[ 2. ,  2. ],
          [ 2. ,  2. ]]])
```

`numina.array.combine.zerocombine` (*arrays, masks, dtype=None, scales=None*)

Combine zero arrays.

#### Parameters

- **arrays** – a list of arrays
- **masks** – a list of mask arrays, True values are masked
- **dtype** – data type of the output
- **scales** –

**Returns** median, variance of the median and number of points stored

### Combination methods in `numina.array.combine`

All these functions return a PyCapsule, that can be passed to `generic_combine()`

`numina.array.combine.mean_method()`

Mean method

`numina.array.combine.median_method()`

Median method

`numina.array.combine.sigmaclip_method([low=0.0[, high=0.0]])`

Sigmaclip method

#### Parameters

- **low** – Number of sigmas to reject under the mean
- **high** – Number of sigmas to reject over the mean

**Raises** `ValueError` if **low** or **high** are negative

`numina.array.combine.quantileclip_method([fclip=0.0])`

Quantile clip method

**Parameters** **fclip** – Fraction of points to reject on both ends

**Raises** `ValueError` if **fclip** is negative or greater than 0.4

```
numina.array.combine.minmax_method([nmin=0[, nmax=0]])
```

Min-max method

#### Parameters

- **nmin** – Number of minimum points to reject
- **nmax** – Number of maximum points to reject

**Raises** `ValueError` if **nmin** or **nmax** are negative

### Extending `generic_combine()`

New combination methods can be implemented and used by `generic_combine()`. The combine function expects a `PyCapsule` object containing a pointer to a C function implementing the combination method.

```
int combine(double *data, double *weights, size_t size, double *out[3], void *func_data)
```

Operate on two arrays, containing **data** and **weights**. The result, its variance and the number of points used in the calculation (useful when there is some kind of rejection) are stored in **out[0]**, **out[1]** and **out[2]**.

#### Parameters

- **data** – a pointer to an array containing the data
- **weights** – a pointer to an array containing weights
- **size** – the size of data and weights
- **out** – an array of pointers to the pixels in the result arrays
- **func\_data** – additional parameters of the function encoded as a void pointer

**Returns** 1 if operation succeeded, 0 in case of error.

If the function uses dynamically allocated data stored in `func_data`, we must also implement a function that deallocates the data once it is used.

```
void destructor_function(PyObject* cobject)
```

#### Parameters

- **cobject** – the object owning dynamically allocated data

### Simple combine method

As an example, I'm going to implement a combination method that returns the minimum of the input arrays. Let's call the method `min_method`

First, we implement the C function. I'm going to use some C++ here (it makes the code very simple).

```
int min_combine(double *data, double *weights, size_t size, double *out[3],
               void *func_data) {

    double* res = std::min_element(data, data + size);

    *out[0] = *res;
    // I'm not going to compute the variance for the minimum
    // but it should go here
    *out[1] = 0.0;
    *out[2] = size;
}
```

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```

return 1;
}

```

A destructor function is not needed in this case as we are not using *func\_data*.

The next step is to build a Python extension. First we need to create a function returning the PyCapsule in C code like this:

```

static PyObject *
py_method_min(PyObject *obj, PyObject *args) {
    if (not PyArg_ParseTuple(args, "")) {
        PyErr_SetString(PyExc_RuntimeError, "invalid parameters");
        return NULL;
    }
    return PyCapsule_New((void*)min_function, "numina.cmethod", NULL);
}

```

The string "numina.cmethod" is the name of the PyCapsule. It cannot be loaded unless it is the name expected by the C code.

The code to load it in a module is like this:

```

static PyMethodDef mymod_methods[] = {
    {"min_combine", (PyCFunction) py_method_min, METH_VARARGS, "Minimum method."},
    ...,
    { NULL, NULL, 0, NULL } /* sentinel */
};

PyMODINIT_FUNC
init_mymodule(void)
{
    PyObject *m;
    m = Py_InitModule("_mymodule", mymod_methods);
}

```

When compiled, this code created a file *\_mymodule.so* that can be loaded by the Python interpreter. This module will contain, among others, a *min\_combine* function.

```

>>> from _mymodule import min_combine
>>> method = min_combine()
...
>>> o = generic_combine(method, arrays)

```

### A combine method with parameters

A combine method with parameters follow a similar approach. Let's say we want to implement a sigma-clipping method. We need to pass the function a *low* and a *high* rejection limits. Both numbers are real numbers greater than zero.

First, the Python function. I'm skipping error checking code here.

```

static PyObject *
py_method_sigmaclip(PyObject *obj, PyObject *args) {
    double low = 0.0;
    double high = 0.0;
}

```

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```

PyObject *cap = NULL;

if (!PyArg_ParseTuple(args, "dd", &low, &high)) {
    PyErr_SetString(PyExc_RuntimeError, "invalid parameters");
    return NULL;
}

cap = PyCapsule_New((void*) my_sigmaclip_function, "numina.cmethod", my_
↳destructor);

/* Allocating space for the two parameters */
/* We use Python memory allocator */
double *funcdata = (double*)PyMem_Malloc(2 * sizeof(double));

funcdata[0] = low;
funcdata[1] = high;
PyCapsule_SetContext(cap, funcdata);
return cap;
}

```

Notice that in this case we construct the `PyCObject` using the same function than in the previous case. The additional data is stored as *Context*.

The deallocator is simply:

```

void my_destructor_function(PyObject* cap) {
    void* cdata = PyCapsule_GetContext(cap);
    PyMem_Free(cdata);
}

```

and the combine function is:

```

int my_sigmaclip_function(double *data, double *weights, size_t size, double *out[3],
    void *func_data) {

    double* fdata = (double*) func_data;
    double slow = *fdata;
    double shigh = *(fdata + 1);

    /* Operations go here */

    return 1;
}

```

Once the module is created and loaded, a sample session would be:

```

>>> from _mymodule import min_combine
>>> method = sigmaclip_combine(3.0, 3.0)
...
>>> o = generic_combine(method, arrays)

```

### numina.array.cosmetics — Array cosmetics

`numina.array.cosmetics.ccdmask` (*flat1*, *flat2=None*, *mask=None*, *lowercut=6.0*, *uppercut=6.0*, *siglev=1.0*, *mode='region'*, *nmed=(7, 7)*, *nsig=(15, 15)*)

Find cosmetic defects in a detector using two flat field images.



Two arrays representing flat fields of different exposure times are required. Cosmetic defects are selected as points that deviate significantly of the expected normal distribution of pixels in the ratio between *flat2* and *flat1*. The median of the ratio is computed and subtracted. Then, the standard deviation is estimated computing the percentiles nearest to the pixel values corresponding to ‘siglev’ in the normal CDF. The standard deviation is then the distance between the pixel values divided by two times *siglev*. The ratio image is then normalized with this standard deviation.

The behavior of the function depends on the value of the parameter *mode*. If the value is ‘region’ (the default), both the median and the sigma are computed in boxes. If the value is ‘full’, these values are computed using the full array.

The size of the boxes in ‘region’ mode is given by *nmed* for the median computation and *nsig* for the standard deviation.

The values in the normalized ratio array above *uppercut* are flagged as hot pixels, and those below ‘-lowercut’ are flagged as dead pixels in the output mask.

#### Parameters

- **flat1** – an array representing a flat illuminated exposure.
- **flat2** – an array representing a flat illuminated exposure.
- **mask** – an integer array representing initial mask.
- **lowercut** – values below this sigma level are flagged as dead pixels.
- **uppercut** – values above this sigma level are flagged as hot pixels.
- **siglev** – level to estimate the standard deviation.
- **mode** – either ‘full’ or ‘region’
- **nmed** – region used to compute the median
- **nsig** – region used to estimate the standard deviation

**Returns** the normalized ratio of the flats, the updated mask and standard deviation

---

**Note:** This function is based on the description of the task `ccdmask` of IRAF

---

#### See also:

***cosmetics()*** Operates much like this function but computes median and sigma in the whole image instead of in boxes

```
numina.array.cosmetics.cosmetics (flat1, flat2=None, mask=None, lowercut=6.0, uppercut=6.0,
                                   siglev=2.0)
```

Find cosmetic defects in a detector using two flat field images.

Two arrays representing flat fields of different exposure times are required. Cosmetic defects are selected as points that deviate significantly of the expected normal distribution of pixels in the ratio between *flat2* and *flat1*.

The median of the ratio array is computed and subtracted to it.

The standard deviation of the distribution of pixels is computed obtaining the percentiles nearest the pixel values corresponding to *nsig* in the normal CDF. The standar deviation is then the distance between the pixel values divided by two times *nsig*. The ratio image is then normalized with this standard deviation.

The values in the ratio above *uppercut* are flagged as hot pixels, and those below ‘-lowercut’ are flagged as dead pixels in the output mask.

#### Parameters

- **flat1** – an array representing a flat illuminated exposure.
- **flat2** – an array representing a flat illuminated exposure.
- **mask** – an integer array representing initial mask.
- **lowercut** – values bellow this sigma level are flagged as dead pixels.
- **uppercut** – values above this sigma level are flagged as hot pixels.
- **siglev** – level to estimate the standard deviation.

**Returns** the updated mask

### `numina.array.fwhm` — FWHM

FWHM calculation

`numina.array.fwhm.compute_fw_at_frac_max_1d_simple(Y, xc, X=None, f=0.5)`  
 Compute the full width at fraction f of the maximum

`numina.array.fwhm.compute_fwhm_1d_simple(Y, xc, X=None)`  
 Compute the FWHM.

### `numina.array.imsurfit` — Image surface fitting

Least squares 2D image fitting to a polynomial.

`numina.array.imsurfit.imsurfit(data, order, output_fit=False)`  
 Fit a bidimensional polynomial to an image.

#### Parameters

- **data** – a bidimensional array
- **order** (*integer*) – order of the polynomial
- **output\_fit** (*bool*) – return the fitted image

**Returns** a tuple with an array with the coefficients of the polynomial terms

```
>>> import numpy
>>> xx, yy = numpy.mgrid[-1:1:100j, -1:1:100j]
>>> z = 456.0 + 0.3 * xx - 0.9* yy
>>> imsurfit(z, order=1) #doctest: +NORMALIZE_WHITESPACE
(array([ 4.56000000e+02,  3.00000000e-01, -9.00000000e-01]),)
```

### `numina.array.interpolation` — Interpolation

A monotonic piecewise cubic interpolator.

**class** `numina.array.interpolation.SteffenInterpolator` (*x*, *y*, *yp\_0=0.0*, *yp\_N=0.0*,  
*extrapolate='raise'*,  
*fill\_value=nan*)

A monotonic piecewise cubic 1-d interpolator.

A monotonic piecewise cubic interpolator based on Steffen, M., *Astronomy & Astrophysics*, 239, 443-450 (1990)

*x* and *y* are arrays of values used to approximate some function  $f: y = f(x)$ . This class returns an object whose call method uses monotonic cubic splines to find the value of new points.

## Parameters

- **x** (*(N,)*, *array\_like*) – A 1-D array of real values, sorted monotonically increasing.
- **y** (*(N,)*, *array\_like*) – A 1-D array of real values.
- **yp\_0** (*float*, *optional*) – The value of the derivative in the first sample.
- **yp\_N** (*float*, *optional*) – The value of the derivative in the last sample.
- **extrapolate** (*str*, *optional*) – Specifies the kind of extrapolation as a string ('extrapolate', 'zeros', 'raise', 'const', 'border') If 'raise' is set, when interpolated values are requested outside of the domain of the input data (x,y), a ValueError is raised. If 'const' is set, 'fill\_value' is returned. If 'zeros' is set, '0' is returned. If 'border' is set, 'y[0]' is returned for values below 'x[0]' and 'y[N-1]' is returned for values above 'x[N-1]' If 'extrapolate' is set, the extreme polynomial are extrapolated outside of their ranges. Default is 'raise'.
- **fill\_value** (*float*, *optional*) – If provided, then this value will be used to fill in for requested points outside of the data range when 'extrapolation' is set to "const". If not provided, then the default is NaN.

## numina.array.mode — Mode

Mode estimators.

`numina.array.mode.mode_half_sample(a, is_sorted=False)`

Estimate the mode using the Half Sample mode.

A method to estimate the mode, as described in D. R. Bickel and R. Frühwirth (contributed equally), "On a fast, robust estimator of the mode: Comparisons to other robust estimators with applications," Computational Statistics and Data Analysis 50, 3500-3530 (2006).

### Example

```
>> import numpy as np >> np.random.seed(1392838) >> a = np.random.normal(1000, 200, size=1000) >>
a[:100] = np.random.normal(2000, 300, size=100) >> b = np.sort(a) >> mode_half_sample(b, is_sorted=True)
1041.9327885039545
```

`numina.array.mode.mode_sex(a)`

Estimate the mode as sextractor

## numina.array.nirproc — nIR preprocessing

`numina.array.nirproc.fowler_array(fowlerdata, ti=0.0, ts=0.0, gain=1.0, ron=1.0, badpixels=None, dtype='float64', saturation=65631, blank=0, normalize=False)`

Loop over the first axis applying Fowler processing.

*fowlerdata* is assumed to be a 3D `numpy.ndarray` containing the result of a nIR observation in Fowler mode (Fowler and Gatley 1991). The shape of the array must be of the form  $2N_p \times M \times N$ , with  $N_p$  being the number of pairs in Fowler mode.

The output signal is just the mean value of the differences between the last  $N_p$  values ( $S_i$ ) and the first  $N_p$  values ( $R_i$ ).

$$S_F = \frac{1}{N_p} \sum_{i=0}^{N_p-1} S_i - R_i$$

If the source has a radiance  $F$ , then the measured signal is equivalent to:

$$S_F = FT_I - FT_S(N_p - 1) = FT_E$$

being  $T_I$  the integration time ( $ti$ ), the time since the first productive read to the last productive read for a given pixel and  $T_S$  the time between samples ( $ts$ ).  $T_E$  is the time between correlated reads  $T_E = T_I - T_S(N_p - 1)$ .

The variance of the signal is the sum of two terms, one for the readout noise:

$$\text{var}(S_{F1}) = \frac{2\sigma_R^2}{N_p}$$

and other for the photon noise:

$$\text{var}(S_{F2}) = FT_E - FT_S \frac{1}{3} \left( N_p - \frac{1}{N_p} \right) = FT_I - FT_S \left( \frac{4}{3} N_p - 1 - \frac{1}{3N_p} \right)$$

### Parameters

- **fowlerdata** – Convertible to a 3D numpy.ndarray with first axis even
- **ti** – Integration time.
- **ts** – Time between samples.
- **gain** – Detector gain.
- **ron** – Detector readout noise in counts.
- **badpixels** – An optional MxN mask of dtype ‘uint8’.
- **dtype** – The dtype of the float outputs.
- **saturation** – The saturation level of the detector.
- **blank** – Invalid values in output are substituted by *blank*.

**Returns** A tuple of (signal, variance of the signal, number of pixels used and badpixel mask).

**Raises** ValueError

```
numina.array.nirproc.ramp_array(rampdata, ti, gain=1.0, ron=1.0, badpixels=None,
                               dtype='float64', saturation=65631, blank=0, nsig=None,
                               normalize=False)
```

Loop over the first axis applying ramp processing.

*rampdata* is assumed to be a 3D numpy.ndarray containing the result of a nIR observation in follow-up-the-ramp mode. The shape of the array must be of the form  $N_s \times M \times N$ , with  $N_s$  being the number of samples.

### Parameters

- **fowlerdata** – Convertible to a 3D numpy.ndarray
- **ti** – Integration time.
- **gain** – Detector gain.
- **ron** – Detector readout noise in counts.
- **badpixels** – An optional MxN mask of dtype ‘uint8’.
- **dtype** – The dtype of the float outputs.
- **saturation** – The saturation level of the detector.
- **blank** – Invalid values in output are substituted by *blank*.

**Returns** A tuple of signal, variance of the signal, number of pixels used and badpixel mask.

**Raises** ValueError

## numina.array.offrot — Offset and Rotation

Fit offset and rotation

numina.array.offrot.**fit\_offset\_and\_rotation**(*coords0*, *coords1*)

Fit a rotation and a traslation between two sets points.

Fit a rotation matrix and a traslation bewtween two matched sets consisting of M N-dimensional points

### Parameters

- **coords0** (*M*, *N*) *array\_like*—
- **coords1** (*M*, *N*) *array\_lke*—

### Returns

- **offset** (*N*, ) *array\_like*
- **rotation** (*N*, *N*) *array\_like*

### Notes

Fit offset and rotation using Kabsch’s algorithm<sup>12</sup>

## numina.array.peaks — Peak finding

## numina.array.recenter — Recenter

Recenter routines

numina.array.recenter.**centering\_centroid**(*data*, *xi*, *yi*, *box*, *nloop=10*, *toldist=0.001*,  
*maxdist=10.0*)

returns x, y, background, status, message

**status is:**

- 0: not recentering
- 1: recentering successful
- 2: maximum distance reached
- 3: not converged

## numina.array.robustfit — Robust fits

Robust fits

numina.array.robustfit.**fit\_theil\_sen**(*x*, *y*)

Compute a robust linear fit using the Theil-Sen method.

See [http://en.wikipedia.org/wiki/Theil%E2%80%93Sen\\_estimator](http://en.wikipedia.org/wiki/Theil%E2%80%93Sen_estimator) for details. This function “pairs up sample points by the rank of their x-coordinates (the point with the smallest coordinate being paired with the first point above the median coordinate, etc.) and computes the median of the slopes of the lines determined by these pairs of points”.

### Parameters

<sup>1</sup> Kabsch algorithm: [https://en.wikipedia.org/wiki/Kabsch\\_algorithm](https://en.wikipedia.org/wiki/Kabsch_algorithm)

<sup>2</sup> Also here: [http://ngihaio.com/?page\\_id=671](http://ngihaio.com/?page_id=671)

- **x** (*array\_like, shape (M,)*) – X coordinate array.
- **y** (*array\_like, shape (M,) or (M,K)*) – Y coordinate array. If the array is two dimensional, each column of the array is independently fitted sharing the same x-coordinates. In this last case, the returned intercepts and slopes are also 1d numpy arrays.

**Returns coef** – Intercept and slope of the linear fit. If y was 2-D, the coefficients in column k of coef represent the linear fit to the data in y’s k-th column.

**Return type** ndarray, shape (2,) or (2, K)

**Raises** ValueError: – If the number of points to fit is < 5

### numina.array.stats —

numina.array.stats.**robust\_std**(*x, debug=False*)

Compute a robust estimator of the standard deviation

See Eq. 3.36 (page 84) in Statistics, Data Mining, and Machine in Astronomy, by Ivezić, Connolly, VanderPlas & Gray

#### Parameters

- **x** (*1d numpy array, float*) – Array of input values which standard deviation is requested.
- **debug** (*bool*) – If True prints computed values

**Returns sigmag** – Robust estimator of the standar deviation

**Return type** float

numina.array.stats.**summary**(*x, rm\_nan=False, debug=False*)

Compute basic statistical parameters.

#### Parameters

- **x** (*1d numpy array, float*) – Input array with values which statistical properties are requested.
- **rm\_nan** (*bool*) – If True, filter out NaN values before computing statistics.
- **debug** (*bool*) – If True prints computed values.

**Returns result** – Number of points, minimum, percentile 25, percentile 50 (median), mean, percentile 75, maximum, standard deviation, robust standard deviation, percentile 15.866 (equivalent to -1 sigma in a normal distribution) and percentile 84.134 (+1 sigma).

**Return type** Python dictionary

### numina.array.trace — Spectrum tracing

### numina.array.wavecalib — Wavelength calibration

Automatic identification of lines and wavelength calibration

```
numina.array.wavecalib.arccalibration.arccalibration(wv_master,          xpos_arc,
                                                    naxis1_arc,          crpix1,
                                                    wv_ini_search, wv_end_search,
                                                    wvmin_useful,          wv-
                                                    max_useful,          er-
                                                    ror_xpos_arc, times_sigma_r,
                                                    frac_triplets_for_sum,
                                                    times_sigma_theil_sen,
                                                    poly_degree_wfit,
                                                    times_sigma_polfilt,
                                                    times_sigma_cook,
                                                    times_sigma_inclusion, ge-
                                                    ometry=None, debugplot=0)
```

Performs arc line identification for arc calibration.

This function is a wrapper of two functions, which are responsible of computing all the relevant information concerning the triplets generated from the master table and the actual identification procedure of the arc lines, respectively.

The separation of those computations in two different functions helps to avoid the repetition of calls to the first function when calibrating several arcs using the same master table.

#### Parameters

- **wv\_master** (*1d numpy array, float*) – Array with wavelengths corresponding to the master table (Angstroms).
- **xpos\_arc** (*1d numpy array, float*) – Location of arc lines (pixels).
- **naxis1\_arc** (*int*) – NAXIS1 for arc spectrum.
- **crpix1** (*float*) – CRPIX1 value to be employed in the wavelength calibration.
- **wv\_ini\_search** (*float*) – Minimum expected wavelength in spectrum.
- **wv\_end\_search** (*float*) – Maximum expected wavelength in spectrum.
- **wvmin\_useful** (*float*) – If not None, this value is used to clip detected lines below it.
- **wvmax\_useful** (*float*) – If not None, this value is used to clip detected lines above it.
- **error\_xpos\_arc** (*float*) – Error in arc line position (pixels).
- **times\_sigma\_r** (*float*) – Times sigma to search for valid line position ratios.
- **frac\_triplets\_for\_sum** (*float*) – Fraction of distances to different triplets to sum when computing the cost function.
- **times\_sigma\_theil\_sen** (*float*) – Number of times the (robust) standard deviation around the linear fit (using the Theil-Sen method) to reject points.
- **poly\_degree\_wfit** (*int*) – Degree for polynomial fit to wavelength calibration.
- **times\_sigma\_polfilt** (*float*) – Number of times the (robust) standard deviation around the polynomial fit to reject points.
- **times\_sigma\_cook** (*float*) – Number of times the standard deviation of Cook’s distances to detect outliers. If zero, this method of outlier detection is ignored.
- **times\_sigma\_inclusion** (*float*) – Number of times the (robust) standard deviation around the polynomial fit to include a new line in the set of identified lines.
- **geometry** (*tuple (4 integers) or None*) – x, y, dx, dy values employed to set the window geometry.

- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed. The valid codes are defined in `numina.array.display.pause_debugplot`.

**Returns** `list_of_wvfeatures` – A list of size equal to the number of identified lines, which elements are instances of the class `WavecalFeature`, containing all the relevant information concerning the line identification.

**Return type** `list` (of `WavecalFeature` instances)

```
numina.array.wavecalib.arccalibration.arccalibration_direct (wv_master,
                                                            ntriplets_master, ratios_master_sorted,
                                                            triplets_master_sorted_list,
                                                            xpos_arc,
                                                            naxis1_arc,      cr-
                                                            pix1, wv_ini_search,
                                                            wv_end_search,
                                                            wvmin_useful=None,
                                                            wv-
                                                            max_useful=None,
                                                            error_xpos_arc=1.0,
                                                            times_sigma_r=3.0,
                                                            frac_triplets_for_sum=0.5,
                                                            times_sigma_theil_sen=10.0,
                                                            poly_degree_wfit=3,
                                                            times_sigma_polfilt=10.0,
                                                            times_sigma_cook=10.0,
                                                            times_sigma_inclusion=5.0,
                                                            geometry=None, de-
                                                            bugplot=0)
```

Performs line identification for arc calibration using line triplets.

This function assumes that a previous call to the function responsible for the computation of information related to the triplets derived from the master table has been previously executed.

#### Parameters

- **wv\_master** (*1d numpy array, float*) – Array with wavelengths corresponding to the master table (Angstroms).
- **ntriplets\_master** (*int*) – Number of triplets built from master table.
- **ratios\_master\_sorted** (*1d numpy array, float*) – Array with values of the relative position of the central line of each triplet, sorted in ascending order.
- **triplets\_master\_sorted\_list** (*list of tuples*) – List with tuples of three numbers, corresponding to the three line indices in the master table. The list is sorted to be in correspondence with `ratios_master_sorted`.
- **xpos\_arc** (*1d numpy array, float*) – Location of arc lines (pixels).
- **naxis1\_arc** (*int*) – NAXIS1 for arc spectrum.
- **crpix1** (*float*) – CRPIX1 value to be employed in the wavelength calibration.
- **wv\_ini\_search** (*float*) – Minimum expected wavelength in spectrum.
- **wv\_end\_search** (*float*) – Maximum expected wavelength in spectrum.
- **wvmin\_useful** (*float or None*) – If not `None`, this value is used to clip detected lines below it.



- **wvmax\_useful** (*float* or *None*) – If not *None*, this value is used to clip detected lines above it.
- **error\_xpos\_arc** (*float*) – Error in arc line position (pixels).
- **times\_sigma\_r** (*float*) – Times sigma to search for valid line position ratios.
- **frac\_triplets\_for\_sum** (*float*) – Fraction of distances to different triplets to sum when computing the cost function.
- **times\_sigma\_theil\_sen** (*float*) – Number of times the (robust) standard deviation around the linear fit (using the Theil-Sen method) to reject points.
- **poly\_degree\_wfit** (*int*) – Degree for polynomial fit to wavelength calibration.
- **times\_sigma\_polfilt** (*float*) – Number of times the (robust) standard deviation around the polynomial fit to reject points.
- **times\_sigma\_cook** (*float*) – Number of times the standard deviation of Cook’s distances to detect outliers. If zero, this method of outlier detection is ignored.
- **times\_sigma\_inclusion** (*float*) – Number of times the (robust) standard deviation around the polynomial fit to include a new line in the set of identified lines.
- **geometry** (*tuple* (4 *integers*) or *None*) – *x*, *y*, *dx*, *dy* values employed to set the window geometry.
- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed. The valid codes are defined in `numina.array.display.pause_debugplot`.

**Returns** **list\_of\_wvfeatures** – A list of size equal to the number of identified lines, which elements are instances of the class `WavecalFeature`, containing all the relevant information concerning the line identification.

**Return type** *list* (of `WavecalFeature` instances)

```
numina.array.wavecalib.arccalibration.fit_list_of_wvfeatures (list_of_wvfeatures,
                                                            naxis1_arc, crpix1,
                                                            poly_degree_wfit,
                                                            weighted=False,
                                                            plot_title=None,
                                                            geometry=None,
                                                            debugplot=0)
```

Fit polynomial to arc calibration `list_of_wvfeatures`.

#### Parameters

- **list\_of\_wvfeatures** (*list* (of `WavecalFeature` instances)) – A list of size equal to the number of identified lines, which elements are instances of the class `WavecalFeature`, containing all the relevant information concerning the line identification.
- **naxis1\_arc** (*int*) – NAXIS1 of arc spectrum.
- **crpix1** (*float*) – CRPIX1 value to be employed in the wavelength calibration.
- **poly\_degree\_wfit** (*int*) – Polynomial degree corresponding to the wavelength calibration function to be fitted.
- **weighted** (*bool*) – Determines whether the polynomial fit is weighted or not, using as weights the values of the cost function obtained in the line identification. Since the weights can be very different, typically weighted fits are not good because, in practice, they totally ignore the points with the smallest weights (which, in the other hand, are useful when handling the borders of the wavelength calibration range).

- **plot\_title** (*string or None*) – Title for residuals plot.
- **geometry** (*tuple (4 integers) or None*) – x, y, dx, dy values employed to set the window geometry.
- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed. The valid codes are defined in `numina.array.display.pause_debugplot`.

**Returns** **solution\_wv** – Instance of class `SolutionArcCalibration`, containing the information concerning the arc lines that have been properly identified. The information about all the lines (including those initially found but at the end discarded) is stored in the list of `WavecalFeature` instances ‘`list_of_wvfeatures`’.

**Return type** `SolutionArcCalibration` instance

```
numina.array.wavecalib.arccalibration.gen_triplets_master(wv_master,      geom-
                                                         etry=None,      debug-
                                                         plot=0)
```

Compute information associated to triplets in master table.

Determine all the possible triplets that can be generated from the array `wv_master`. In addition, the relative position of the central line of each triplet is also computed.

**Parameters**

- **wv\_master** (*1d numpy array, float*) – Array with wavelengths corresponding to the master table (Angstroms).
- **geometry** (*tuple (4 integers) or None*) – x, y, dx, dy values employed to set the window geometry.
- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed. The valid codes are defined in `numina.array.display.pause_debugplot`.

**Returns**

- **ntriplets\_master** (*int*) – Number of triplets built from master table.
- **ratios\_master\_sorted** (*1d numpy array, float*) – Array with values of the relative position of the central line of each triplet, sorted in ascending order.
- **triplets\_master\_sorted\_list** (*list of tuples*) – List with tuples of three numbers, corresponding to the three line indices in the master table. The list is sorted to be in correspondence with `ratios_master_sorted`.

```
numina.array.wavecalib.arccalibration.match_wv_arrays(wv_master,
                                                       wv_expected_all_peaks,
                                                       delta_wv_max)
```

Match two lists with wavelengths.

Assign individual wavelengths from `wv_master` to each expected wavelength when the latter is within the maximum allowed range.

**Parameters**

- **wv\_master** (*numpy array*) – Array containing the master wavelengths.
- **wv\_expected\_all\_peaks** (*numpy array*) – Array containing the expected wavelengths (computed, for example, from an approximate polynomial calibration applied to the location of the line peaks).
- **delta\_wv\_max** (*float*) – Maximum distance to accept that the master wavelength corresponds to the expected wavelength.

**Returns** **wv\_verified\_all\_peaks** – Verified wavelengths from master list.

**Return type** numpy array

```
numina.array.wavecalib.arccalibration.refine_arccalibration(sp, poly_initial,
                                                         wv_master,
                                                         poldeg, nrepeat=3,
                                                         ntimes_match_wv=2,
                                                         nwin-
                                                         width_initial=7,
                                                         nwin-
                                                         width_refined=5,
                                                         times_sigma_reject=5,
                                                         interactive=False,
                                                         threshold=0, plot-
                                                         title=None, dec-
                                                         imal_places=4,
                                                         ylogscale=False,
                                                         geometry=None,
                                                         pdf=None, debug-
                                                         plot=0)
```

Refine wavelength calibration using an initial polynomial.

#### Parameters

- **sp** (*numpy array*) – 1D array of length NAXIS1 containing the input spectrum.
- **poly\_initial** (*Polynomial instance*) – Initial wavelength calibration polynomial, providing the wavelength as a function of pixel number (running from 1 to NAXIS1).
- **wv\_master** (*numpy array*) – Array containing the master list of arc line wavelengths.
- **poldeg** (*int*) – Polynomial degree of refined wavelength calibration. Note that this degree can be different from the polynomial degree of `poly_initial`.
- **nrepeat** (*int*) – Number of times lines are iteratively included in the initial fit.
- **ntimes\_match\_wv** (*int*) – Number of pixels around each line peak where the expected wavelength must match the tabulated wavelength in the master list.
- **nwinwidth\_initial** (*int*) – Initial window width to search for line peaks in spectrum.
- **nwinwidth\_refined** (*int*) – Window width to refine line peak location.
- **times\_sigma\_reject** (*float*) – Times sigma to reject points in the fit.
- **interactive** (*bool*) – If True, the function allows the user to modify the fit interactively.
- **threshold** (*float*) – Minimum signal in the peaks.
- **plottitle** (*string or None*) – Plot title.
- **decimal\_places** (*int*) – Number of decimal places to be employed when displaying relevant fitted parameters.
- **ylogscale** (*bool*) – If True, the spectrum is displayed in logarithmic units. Note that this is only employed for display purposes. The line peaks are found in the original spectrum.
- **geometry** (*tuple (4 integers) or None*) – x, y, dx, dy values employed to set the window geometry.
- **pdf** (*PdfFile object or None*) – If not None, output is sent to PDF file.
- **debugplot** (*int*) – Debugging level for messages and plots. For details see ‘numina.array.display.pause\_debugplot.py’.

**Returns**

- **poly\_refined** (*Polynomial instance*) – Refined wavelength calibration polynomial.
- **yres\_summary** (*dictionary*) – Statistical summary of the residuals.

`numina.array.wavecalib.arccalibration.select_data_for_fit` (*list\_of\_wvfeatures*)

Select information from valid arc lines to facilitate posterior fits.

**Parameters** `list_of_wvfeatures` (*list (of WavecalFeature instances)*) – A list of size equal to the number of identified lines, which elements are instances of the class `WavecalFeature`, containing all the relevant information concerning the line identification.

**Returns**

- **nfit** (*int*) – Number of valid points for posterior fits.
- **ifit** (*list of int*) – List of indices corresponding to the arc lines which coordinates are going to be employed in the posterior fits.
- **xfit** (*1d numpy array*) – X coordinate of points for posterior fits.
- **yfit** (*1d numpy array*) – Y coordinate of points for posterior fits.
- **wfit** (*1d numpy array*) – Cost function of points for posterior fits. The inverse of these values can be employed for weighted fits.

`numina.array.wavecalib.peaks_spectrum.find_peaks_spectrum` (*sx, nwinwidth, threshold=0, debugplot=0*)

Find peaks in array.

The algorithm imposes that the signal at both sides of the peak decreases monotonically.

**Parameters**

- **sx** (*1d numpy array, floats*) – Input array.
- **nwinwidth** (*int*) – Width of the window where each peak must be found.
- **threshold** (*float*) – Minimum signal in the peaks.
- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed: 00 : no debug, no plots 01 : no debug, plots without pauses 02 : no debug, plots with pauses 10 : debug, no plots 11 : debug, plots without pauses 12 : debug, plots with pauses

**Returns** `ixpeaks` – Peak locations, in array coordinates (integers).

**Return type** 1d numpy array, int

`numina.array.wavecalib.peaks_spectrum.refine_peaks_spectrum` (*sx, ixpeaks, nwinwidth, method=None, geometry=None, debugplot=0*)

Refine line peaks in spectrum.

**Parameters**

- **sx** (*1d numpy array, floats*) – Input array.
- **ixpeaks** (*1d numpy array, int*) – Initial peak locations, in array coordinates (integers). These values can be the output from the function `find_peaks_spectrum()`.
- **nwinwidth** (*int*) – Width of the window where each peak must be refined.
- **method** (*string*) – “poly2” : fit to a 2nd order polynomial “gaussian” : fit to a Gaussian

- **geometry** (*tuple (4 integers) or None*) – x, y, dx, dy values employed to set the window geometry.
- **debugplot** (*int*) – Determines whether intermediate computations and/or plots are displayed: 00 : no debug, no plots 01 : no debug, plots without pauses 02 : no debug, plots with pauses 10 : debug, no plots 11 : debug, plots without pauses 12 : debug, plots with pauses

#### Returns

- **fxpeaks** (*1d numpy array, float*) – Refined peak locations, in array coordinates.
- **sxpeaks** (*1d numpy array, float*) – When fitting Gaussians, this array stores the fitted line widths (sigma). Otherwise, this array returns zeros.

Store the solution of a wavelength calibration

**class** numina.array.wavecalib.solutionarc.**CrLinear** (*crpix, crval, crmin, crmax, cdelt*)  
Store information concerning the linear wavelength calibration.

#### Parameters

- **crpix** (*float*) – CRPIX1 value employed in the linear wavelength calibration.
- **crval** (*float*) – CRVAL1 value corresponding to the linear wavelength calibration.
- **crmin** (*float*) – CRVAL value at pixel number 1 corresponding to the linear wavelength calibration.
- **crmax** (*float*) – CRVAL value at pixel number NAXIS1 corresponding to the linear wavelength calibration.
- **cdelt** (*float*) – CDELTA1 value corresponding to the linear wavelength calibration.

Identical to parameters.

**class** numina.array.wavecalib.solutionarc.**SolutionArcCalibration** (*features, coeff, residual\_std, cr\_linear*)

Auxiliary class to store the arc calibration solution.

Note that this class only stores the information concerning the arc lines that have been properly identified. The information about all the lines (including those initially found but at the end discarded) is stored in the list of WavecalFeature instances.

#### Parameters

- **features** (*list (of WavecalFeature instances)*) – A list of size equal to the number of identified lines, which elements are instances of the class WavecalFeature, containing all the relevant information concerning the line identification.
- **coeff** (*1d numpy array (float)*) – Coefficients of the wavelength calibration polynomial.
- **residual\_std** (*float*) – Residual standard deviation of the fit.
- **cr\_linear** (*instance of CrLinear*) – Object containing the linear approximation parameters crpix, crval, cdelt, crmin and crmax.

Identical to parameters.

**update\_features** (*poly*)

Evaluate wavelength at xpos using the provided polynomial.

**class** numina.array.wavecalib.solutionarc.**WavecalFeature** (*line\_ok, category, lineid, funcost, xpos, ypos=0.0, peak=0.0, fwhm=0.0, reference=0.0, wavelength=0.0*)

Store information concerning a particular line identification.

**Parameters**

- **line\_ok** (*bool*) – True if the line has been properly identified.
- **category** (*char*) – Line identification type (A, B, C, D, E, R, T, P, K, I, X). See documentation embedded within the arccalibration\_direct function for details.
- **lineid** (*int*) – Number of identified line within the master list.
- **xpos** (*float*) – Pixel x-coordinate of the peak of the line.
- **ypos** (*float*) – Pixel y-coordinate of the peak of the line.
- **peak** (*float*) – Flux of the peak of the line.
- **fwhm** (*float*) – FWHM of the line.
- **reference** (*float*) – Wavelength of the identified line in the master list.
- **wavelength** (*float*) – Wavelength of the identified line estimated from the wavelength calibration polynomial.
- **funcost** (*float*) – Cost function corresponding to each identified arc line.

**Identical to parameters.**

**numina.array.utils —**

Utility routines

numina.array.utils.**coor\_to\_pix\_1d** (*w*)

Return the pixel where a coordinate is located.

numina.array.utils.**expand\_region** (*tuple\_of\_s, a, b, start=0, stop=None*)

Apply expand\_slice on a tuple of slices

numina.array.utils.**expand\_slice** (*s, a, b, start=0, stop=None*)

Expand a slice on the start/stop limits

numina.array.utils.**image\_box** (*center, shape, box*)

Create a region of size box, around a center in a image of shape.

numina.array.utils.**slice\_create** (*center, block, start=0, stop=None*)

Return an slice with a symmetric region around center.

### 3.1.2 numina.core — Core classes for Pipelines

**numina.core.dataholders — Dataholders**

Recipe requirements

```
class numina.core.dataholders.Parameter (value, description, destination=None, optional=True, choices=None, validation=True, validator=None, accept_scalar=False, as_list=False, nelem=None)
```

The Recipe requires a plain Python type.

#### Parameters

- **value** (*plain python type*) – Default value of the parameter, the requested type is inferred from the type of value.
- **description** (*str*) – Description of the parameter. The value is used by *numina show-recipes* to provide human-readable documentation.
- **destination** (*str, optional*) – Name of the field in the RecipeInput object. Overrides the value provided by the name of the Parameter variable
- **optional** (*bool, optional*) – If *False*, the builder of the RecipeInput must provide a value for this Parameter. If *True* (default), the builder can skip this Parameter and then the default in *value* is used.
- **choices** (*list of plain python type, optional*) – The possible values of the inputs. Any other value will raise an exception
- **validator** (*callable, optional*) – A custom validator for inputs
- **accept\_scalar** (*bool, optional*) – If *True*, when *value* is a list, scalar value inputs are converted to list. If *False* (default), scalar values will raise an exception if *value* is a list
- **as\_list** (*bool, optional*:) – If *True*, consider the internal type a list even if *value* is scalar Default is *False*
- **nelem** (*str or int, optional*:) – If *nelem* is '\*', the list can contain any number of objects. If is '+', the list must contain at least 1 element. With a number, the list must contain that number of elements.

**convert** (*val*)

Convert input values to type values.

**validate** (*val*)

Validate values according to the requirement

```
class numina.core.dataholders.Product (ptype, description="", validation=True, destination=None, optional=False, default=None, choices=None)
```

Product holder for RecipeResult.

Deprecated since version 0.16: *Product* is replaced by *Result*. It will be removed in 1.0

```
class numina.core.dataholders.Requirement (rtype, description, destination=None, optional=False, default=None, choices=None, validation=True, query_opts=None)
```

Requirement holder for RecipeInput.

#### Parameters

- **rtype** (*DataType* or *Type[DataType]*) – Object or class representing the type of the requirement, it must be a subclass of *DataType*
- **description** (*str*) – Description of the Requirement. The value is used by *numina show-recipes* to provide human-readable documentation.
- **destination** (*str, optional*) – Name of the field in the RecipeInput object. Overrides the value provided by the name of the Requirement variable

- **optional** (*bool, optional*) – If *False*, the builder of the `RecipeInput` must provide a value for this Parameter. If *True* (default), the builder can skip this Parameter and then the default in *value* is used.
- **default** (*optional*) – The value provided by the Requirement if the `RecipeInput` builder does not provide one.
- **choices** (*list of values, optional*) – The possible values of the inputs. Any other value will raise an exception

```
class numina.core.dataholders.Result (ptype, description="", validation=True, destination=None, optional=False, default=None, choices=None)
```

Result holder for `RecipeResult`.

### numina.core.metaclass — Metaclasses

Base metaclasses

```
class numina.core.metaclass.RecipeInputType  
    Metaclass for RecipeInput.
```

```
class numina.core.metaclass.RecipeResultType  
    Metaclass for RecipeResult.
```

```
class numina.core.metaclass.StoreType  
    Metaclass for storing members.
```

### numina.core.metarecipes — Meta class for recipes

Metaclasses for Recipes.

```
class numina.core.metarecipes.RecipeType  
    Metaclass for Recipe.
```

```
numina.core.metarecipes.generate_docs (klass)  
    Add documentation to generated classes
```

### numina.core.oresult — Observation Result

Results of the Observing Blocks

```
class numina.core.oresult.ObservationResult (instrument='UNKNOWN', mode='UNKNOWN')
```

The result of a observing block.

```
get_sample_frame ()  
    Return first available frame in observation result
```

```
numina.core.oresult.dataframe_from_list (values)  
    Build a DataFrame object from a list.
```

```
numina.core.oresult.oblock_from_dict (values)  
    Build a ObservingBlock object from a dictionary.
```

```
numina.core.oresult.obsres_from_dict (values)  
    Build a ObservationResult object from a dictionary.
```



**numina.core.pipeline — Pipeline classes**

DRP related classes

```
class numina.core.pipeline.InstrumentDRP (name, configurations, modes, pipelines,
                                         products=None, datamodel=None, ver-
                                         sion='undefined')
```

Description of an Instrument Data Reduction Pipeline

**Parameters**

- **name** (*str*) – Name of the instrument
- **configurations** (*dict of InstrumentConfiguration*) –
- **modes** (*dict of ObservingModes*) –
- **pipeline** (*dict of Pipeline*) –

```
get_recipe_object (mode_name, pipeline_name='default')
    Build a recipe object from a given mode name
```

```
iterate_mode_provides (modes, pipeline)
    Return the mode that provides a given product
```

```
query_provides (product, pipeline='default', search=False)
    Return the mode that provides a given product
```

```
search_mode_provides (product, pipeline='default')
    Search the mode that provides a given product
```

```
select_configuration_old (obresult)
    Select instrument configuration based on OB
```

```
select_profile (obresult)
    Select instrument profile based on OB
```

```
select_profile_image (img)
    Select instrument profile based on FITS
```

```
class numina.core.pipeline.ObservingMode
    Observing modes of an Instrument.
```

```
class numina.core.pipeline.Pipeline (instrument, name, recipes, version=1, products=None,
                                       provides=None)
```

Base class for pipelines.

```
depsolve ()
    Load all recipes to search for products
```

```
get_recipe_object (mode)
    Load recipe object, according to observing mode
```

```
load_product_class (mode)
    Load recipe object, according to observing mode
```

```
load_product_object (name)
    Load product object, according to name
```

```
load_recipe_object (mode)
    Load recipe object, according to observing mode
```

```
provides (mode_label)
    Return the ProductEntry for some mode
```

**query\_recipe** (*mode*)  
Recursive query of all calibrations required by a mode

**who\_provides** (*product\_label*)  
Return the ProductEntry for some requirement

### **numina.core.pipelineload** — Build pipelines from files

Build a LoadableDRP from a yaml file

**numina.core.pipelineload.check\_section** (*node, section, keys=None*)  
Validate keys in a section

**numina.core.pipelineload.drp\_load** (*package, resource, confclass=None*)  
Load the DRPS from a resource file.

**numina.core.pipelineload.drp\_load\_data** (*package, data, confclass=None*)  
Load the DRPS from data.

**numina.core.pipelineload.load\_link** (*node*)  
Build a ResultOf query from dict

**numina.core.pipelineload.load\_mode** (*node*)  
Load one observing mode

**numina.core.pipelineload.load\_mode\_builder** (*obs\_mode, node*)  
Load observing mode OB builder

**numina.core.pipelineload.load\_mode\_tagger** (*obs\_mode, node*)  
Load observing mode OB tagger

**numina.core.pipelineload.load\_mode\_validator** (*obs\_mode, node*)  
Load observing mode validator

**numina.core.pipelineload.load\_modes** (*node*)  
Load all observing modes

### **numina.core.recipeinout** — Recipe input and output

Recipe inputs and outputs

**class** **numina.core.recipeinout.RecipeInput** (*\*args, \*\*kws*)  
RecipeInput base class

**class** **numina.core.recipeinout.RecipeResult** (*\*args, \*\*kws*)

**class** **numina.core.recipeinout.RecipeResultBase** (*\*args, \*\*kws*)  
The result of a Recipe.

**class** **numina.core.recipeinout.define\_input** (*input\_class*)  
Recipe decorator.

**numina.core.recipeinout.define\_requirements**  
alias of *numina.core.recipeinout.define\_input*

**class** **numina.core.recipeinout.define\_result** (*resultClass*)  
Recipe decorator.

## numina.core.recipes — Base class for Recipes

Basic tools and classes used to generate recipe modules.

A recipe is a class that complies with the *reduction recipe API*:

- The class must derive from `numina.core.BaseRecipe`.

**class** `numina.core.recipes.BaseRecipe` (*\*args, \*\*kwargs*)

Base class for all instrument recipes

**Parameters** `intermediate_results` (*bool, optional*) – If True, save intermediate results of the Recipe

**obresult**

**Type** `ObservationResult`, requirement

**logger**

recipe logger

**class** `RecipeInput` (*\*args, \*\*kwargs*)

RecipeInput base class

**class** `RecipeResult` (*\*args, \*\*kwargs*)

**build\_recipe\_input** (*ob, dal*)

Build a RecipeInput object.

**classmethod** `create_input` (*\*args, \*\*kwargs*)

Pass the result arguments to the RecipeInput constructor

**classmethod** `create_result` (*\*args, \*\*kwargs*)

Pass the result arguments to the RecipeResult constructor

**run\_qc** (*recipe\_input, recipe\_result*)

Run Quality Control checks.

**save\_intermediate\_array** (*array, name*)

Save intermediate array object as FITS.

**save\_intermediate\_img** (*img, name*)

Save intermediate FITS objects.

**set\_base\_headers** (*hdr*)

Set metadata in FITS headers.

**validate\_input** (*recipe\_input*)

“Validate the input of the recipe

**validate\_result** (*recipe\_result*)

Validate the result of the recipe

`numina.core.recipes.timeit` (*method*)

Decorator to measure the time used by the recipe

## numina.core.requirements — Recipe requirements

Recipe requirement holders

**class** `numina.core.requirements.ObservationResultRequirement` (*query\_opts=None*)

The Recipe requires the result of an observation.

### numina.core.taggers — Extract information from OBs

Function to retrieve tags from Observation results.

`numina.core.taggers.get_tags_from_full_ob` (*ob*, *reqtags=None*)

#### Parameters

- (**ObservationResult**) (*ob*) –
- (**dict**) (*reqtags*) –

#### Returns

**Return type** A dictionary

### numina.core.types — Types for Recipe IO

Deprecated since version 0.17: Use `numina.types`

### numina.core.utils — Utilities

Recipes for system checks.

**class** `numina.core.utils.AlwaysFailRecipe` (*\*args*, *\*\*kwargs*)

A Recipe that always fails.

**class** `RecipeInput` (*\*args*, *\*\*kwargs*)

RecipeInput base class

**class** `RecipeResult` (*\*args*, *\*\*kwargs*)

**class** `numina.core.utils.AlwaysSuccessRecipe` (*\*args*, *\*\*kwargs*)

A Recipe that always successes.

**class** `RecipeInput` (*\*args*, *\*\*kwargs*)

RecipeInput base class

**class** `RecipeResult` (*\*args*, *\*\*kwargs*)

**class** `numina.core.utils.Combine` (*\*args*, *\*\*kwargs*)

**class** `CombineInput` (*\*args*, *\*\*kwargs*)

CombineInput documentation.

#### field

Extract field of previous result

**Type** `str`, requirement, optional, default=image

#### method

Method of combination

**Type** `str`, requirement, optional, default=mean

#### method\_kwargs

Arguments passed to the combination method

**Type** `dict`, requirement, optional

#### obresult

Observation Result

**Type** `ObservationResultType`, requirement

```
class CombineResult (*args, **kws)
    CombineResult documentation.
```

```
    result
        Type DataFrameType, product
```

```
RecipeInput
    alias of numina.core.metaclass.CombineInput
```

```
RecipeResult
    alias of numina.core.metaclass.CombineResult
```

```
build_recipe_input (obsres, dal)
    Build a RecipeInput object.
```

```
class numina.core.utils.OBSuccessRecipe (*args, **kwargs)
    A Recipe that always successes, it requires an OB
```

```
    class OBSuccessRecipeInput (*args, **kws)
        OBSuccessRecipeInput documentation.
```

```
    obresult
        Observation Result
        Type ObservationResultType, requirement
```

```
    RecipeInput
        alias of numina.core.metaclass.OBSuccessRecipeInput
```

```
    class RecipeResult (*args, **kws)
```

### numina.core.validator — Input output validation

Validator decorator

```
numina.core.validator.as_list (callable)
    Convert a scalar validator in a list validator
```

```
numina.core.validator.only_positive (value)
    Validation error is value is negative
```

```
numina.core.validator.range_validator (minval=None, maxval=None)
    Generates a function that validates that a number is within range
```

#### Parameters

- **minval** (*numeric, optional*) – Values strictly lesser than *minval* are rejected
- **maxval** (*numeric, optional*) – Values strictly greater than *maxval* are rejected

#### Returns

- A function that returns values if are in the range and raises
- *ValidationError* is the values are outside the range

```
numina.core.validator.validate (method)
    Decorate run method, inputs and outputs are validated
```

### 3.1.3 numina.dal — Calibration fetching

DAL base class

**class** numina.dal.daliface.**DALInterface**

Abstract Base Class for DAL

DAL base classes

**class** numina.dal.absdal.**AbsDAL**

**class** numina.dal.absdal.**AbsDrpDAL** (*drps, \*args, \*\*kwargs*)

DAL for dictionary-based database of products.

**class** numina.dal.dictdal.**BaseDictDAL** (*drps, ob\_table, prod\_table, req\_table, extra\_data=None, components=None*)

A dictionary based DAL

**obsres\_from\_oblock\_id** (*obsid, as\_mode=None, configuration=None*)

” Override instrument configuration if configuration is not None

**search\_prod\_obsid** (*ins, obsid, pipeline*)

Returns the first coincidence...

**search\_prod\_type\_tags** (*tipo, ins, tags, pipeline*)

Returns the first coincidence...

**class** numina.dal.dictdal.**BaseHybridDAL** (*drps, obtable, base, extra\_data=None, basedir=None, components=None*)

**class** numina.dal.dictdal.**Dict2DAL** (*drps, obtable, base, extra\_data=None, components=None*)

**class** numina.dal.dictdal.**DictDAL** (*drps, base*)

**class** numina.dal.dictdal.**HybridDAL** (*drps, obtable, base, extra\_data=None, components=None, basedir=None*)

A DAL that can read files from directory structure

DAL for file-based database of products.

DAL Mock class

**class** numina.dal.mockdal.**MockDAL**

**search\_prod\_type\_tags** (*ins, type, tags, pipeline*)

Returns the first coincidence...

**class** numina.dal.stored.**StoredParameter** (*content*)

A parameter returned from the DAL

**class** numina.dal.stored.**StoredProduct** (*id, content, tags, \*\*kwds*)

A product returned from the DAL

**class** numina.dal.stored.**StoredResult**

Recover the RecipeResult values stored in the Backend

DAL for dictionary-based database of products.

**class** numina.dal.backend.**Backend** (*drps, base, extra\_data=None, basedir=None, components=None, filename=None*)

Utilities for DAL

numina.dal.utils.**tags\_are\_valid** (*subset, superset*)

Validate tags

### 3.1.4 numina.datamodel — Model of data

**class** numina.datamodel.**DataModel** (*name='UNKNOWN', mappings=None*)  
 Model of the Data being processed

#### Parameters

- **name** (*str*) – Name of the DataModel
- **mappings** (*dict, optional*) – A dictionary of str keys and values convertibles to FITSKeyExtractor

**do\_sky\_correction** (*img*)  
 Perform sky correction

**gather\_info** (*dframe*)  
 Obtain a summary of information about the image.

**gather\_info\_dframe** (*img*)  
 Obtain a summary of information about the image.

**gather\_info\_hdu** (*hdulist*)  
 Obtain a summary of information about the image.

**get\_darktime** (*img*)  
 Obtain DARKTIME

**get\_data** (*img*)  
 Obtain the primary data from the image

**get\_exptime** (*img*)  
 Obtain EXPTIME

**get\_header** (*img*)  
 Obtain the primary header from the image

**get\_imgid** (*img*)  
 Obtain a unique identifier of the image.

**Parameters** *img* (*astropy.io.fits.HDUList*) –

**Returns** Identification of the image

**Return type** *str*

**get\_quality\_control** (*img*)  
 Obtain quality control flag from the image.

**get\_variance** (*img*)  
 Obtain the variance of the primary data from the image

**class** numina.datamodel.**FITSKeyExtractor** (*values*)  
 Extract values from FITS images

### 3.1.5 numina.drps — Loading DRPS

DRP system wide initialization

numina.drps.**get\_system\_drps** ()  
 Load all compatible DRPs in the system

DRP storage class

**class** numina.drps.drpbased.DrpBase

Store DRPs, base class

**class** numina.drps.drpbased.DrpGeneric (*drps=None*)

Store DRPs in a dictionary

**query\_all** ()

Return all available DRPs in internal storage

**query\_by\_name** (*name*)

Query DRPs in internal storage by name

DRP system-wide loader

**class** numina.drps.drpsystem.DrpSystem (*entry\_point='numina.pipeline.1'*)

Load DRPs from the system.

**classmethod** **iload** (*entry\_point='numina.pipeline.1'*)

Load all available DRPs in 'entry\_point'.

**load** ()

Load all available DRPs in 'entry\_point'.

**classmethod** **load\_drp** (*name, entry\_point='numina.pipeline.1'*)

Load all available DRPs in 'entry\_point'.

### 3.1.6 numina.exceptions — Numina exceptions

Exceptions for the numina package.

**exception** numina.exceptions.DetectorElapseError

Error in the clocking of a Detector.

**exception** numina.exceptions.DetectorReadoutError

Error in the readout of a Detector.

**exception** numina.exceptions.Error

Base class for exceptions in the numina package.

**exception** numina.exceptions.NoResultFound

No result found in a DAL query.

numina.exceptions.NoResultFoundOrig

alias of *numina.exceptions.NoResultFound*

**exception** numina.exceptions.RecipeError

A non recoverable problem during recipe execution.

**exception** numina.exceptions.ValidationError

Error during validation of Recipe inputs and outputs.

### 3.1.7 numina.frame — Frame manipulation

FITS header schema and validation.

This module is a simplification of the FITS Schema defined by Erik Bray here: [http://embray.github.io/PyFITS/schema/users\\_guide/users\\_schema.html](http://embray.github.io/PyFITS/schema/users_guide/users_schema.html)

If this schema implementation reaches pyfits/astropy stable, we will use it instead of ours, with schema definitions being the same.



**class** `numina.frame.schema.Schema` (*sc*)  
A FITS schema

**exception** `numina.frame.schema.SchemaDefinitionError`  
Exception raised when a FITS Schema definition is not valid.

**class** `numina.frame.schema.SchemaKeyword` (*name*, *mandatory=False*, *valid=True*,  
*value=None*)  
A keyword in the schema

**exception** `numina.frame.schema.SchemaValidationError`  
Exception raised when a Schema does not validate a FITS header.

### 3.1.8 `numina.logger` — Logging utilities

Extra logging handlers for the numina logging system.

**class** `numina.logger.FITSHistoryHandler` (*header*)  
Logging handler using HISTORY FITS cards

**emit** (*record*)  
Do whatever it takes to actually log the specified logging record.

This version is intended to be implemented by subclasses and so raises a `NotImplementedError`.

`numina.logger.log_to_history` (*logger*, *name*)  
Decorate function, adding a logger handler stored in FITS.

### 3.1.9 `numina.modeling` — Model for fitting

**class** `numina.modeling.enclosed.EnclosedGaussian` (*amplitude*, *stddev*, *\*\*kwargs*)  
Enclosed Gaussian model

**static evaluate** (*x*, *amplitude*, *stddev*)  
Evaluate the model on some input variables.

**class** `numina.modeling.gaussbox.GaussBox` (*amplitude=1.0*, *mean=0.0*, *stddev=1.0*, *hpix=0.5*,  
*\*\*kwargs*)

Model for fitting a 1D Gaussina convolved with a square

**static evaluate** (*x*, *amplitude*, *mean*, *stddev*, *hpix*)  
Evaluate the model on some input variables.

`numina.modeling.gaussbox.gauss_box_model` (*x*, *amplitude=1.0*, *mean=0.0*, *stddev=1.0*,  
*hpix=0.5*)

Integrate a Gaussian profile.

`numina.modeling.gaussbox.gauss_box_model_deriv` (*x*, *amplitude=1.0*, *mean=0.0*, *std-*  
*dev=1.0*, *hpix=0.5*)

Derivative of the integral of a Gaussian profile.

`numina.modeling.sumofgauss.sum_of_gaussian_factory` (*N*)  
Return a model of the sum of N Gaussians and a constant background.

### 3.1.10 `numina.core.qc` — Quality Control for Numina

This module defines functions and classes which implement quality assesment for Numina-based applications.

Deprecated since version 0.17: Use `numina.types.qc`

## QA Levels

The numeric values of the QC levels are given in this table.

Level	Numeric value
GOOD	100
FAIR	90
BAD	70

### 3.1.11 `numina.store` — Serialization of products

### 3.1.12 `numina.treedict` — Recursive dictionary

An implementation of hierarchical dictionary.

### 3.1.13 `numina.types` — Data types

**class** `numina.types.base.DataTypeBase` (*\*args, \*\*kwargs*)

Base class for input/output types of recipes.

**static** `create_db_info()`

Create metadata structure

**extract\_db\_info** (*obj, db\_info\_keys*)

Extract metadata from serialized file

**classmethod** `isproduct()`

Check if the `DataType` is the product of a `Recipe`

**name** ()

Unique name of the datatype

**update\_meta\_info** ()

Extract my metadata

**class** `numina.types.array.ArrayNTType` (*dimensions, default=None*)

**class** `numina.types.array.ArrayType` (*default=None*)

A type of array.

**convert** (*obj*)

Basic conversion to internal type

This method is intended to be redefined by subclasses

Basic Data Products

**class** `numina.types.dataframe.DataFrame` (*frame=None, filename=None*)

A handle to a image in disk or in memory.

**class** `numina.types.datatype.AnyType`

Type representing anything

**convert** (*obj*)

Basic conversion to internal type

This method is intended to be redefined by subclasses

**validate** (*obj*)  
 Validate convertibility to internal representation

**Returns** True if ‘obj’ matches the data type

**Return type** bool

**Raises** `ValidationError` – If the validation fails

**class** `numina.types.datatype.AutoDataType` (*\*args, \*\*kwargs*)  
 Data type for types that are its own python type

**class** `numina.types.datatype.DataType` (*ptype, node\_type=None, default=None, \*\*kwds*)  
 Base class for input/output types of recipes.

**convert** (*obj*)  
 Basic conversion to internal type

This method is intended to be redefined by subclasses

**convert\_in** (*obj*)  
 Basic conversion to internal type of inputs.

This method is intended to be redefined by subclasses

**convert\_out** (*obj*)  
 Basic conversion to internal type of outputs.

This method is intended to be redefined by subclasses

**validate** (*obj*)  
 Validate convertibility to internal representation

**Returns** True if ‘obj’ matches the data type

**Return type** bool

**Raises** `ValidationError` – If the validation fails

**class** `numina.types.datatype.ListOfType` (*ref, default=None, index=0, nmin=None, nmax=None, accept\_scalar=False, multi\_query=None*)  
 Data type for lists of other types.

**convert** (*obj*)  
 Basic conversion to internal type

This method is intended to be redefined by subclasses

**validate** (*obj*)  
 Validate convertibility to internal representation

**Returns** True if ‘obj’ matches the data type

**Return type** bool

**Raises** `ValidationError` – If the validation fails

**class** `numina.types.datatype.NullType`  
 Data type for None.

**convert** (*obj*)  
 Basic conversion to internal type

This method is intended to be redefined by subclasses

**validate** (*obj*)

Validate convertibility to internal representation

**Returns** True if ‘obj’ matches the data type

**Return type** bool

**Raises** `ValidationError` – If the validation fails

**class** `numina.types.datatype.PlainPythonType` (*ref=None, validator=None*)

Data type for Python basic types.

**convert** (*obj*)

Basic conversion to internal type

This method is intended to be redefined by subclasses

**class** `numina.types.frame.DataFrameType` (*datamodel=None*)

A type of DataFrame.

**convert** (*obj*)

Convert

**extract\_db\_info** (*obj, keys*)

Extract tags from serialized file

**validate** (*value*)

**class** `numina.types.linescatalog.DataProductType` (*ptype, default=None*)

**class** `numina.types.linescatalog.LinesCatalog`

**extract\_db\_info** (*obj, keys*)

Extract metadata from serialized file

**class** `numina.types.multitype.MultiType` (*\*args*)

**isproduct** ()

Check if the DataType is the product of a Recipe

**validate** (*obj*)

Validate convertibility to internal representation

**Returns** True if ‘obj’ matches the data type

**Return type** bool

**Raises** `ValidationError` – If the validation fails

Quality control for Numina-based applications.

**class** `numina.types.qc.QC`

An enumeration.

## QA Levels

The numeric values of the QC levels are given in this table.

Level	Numeric value
GOOD	100
FAIR	90
BAD	70

**class** `numina.types.structured.BaseStructuredCalibration` (*instrument='unknown', datamodel=None*)

Base class for structured calibration data

**Parameters** `instrument` (*str*) – Instrument name

**tags**

dictionary of selection fields

**Type** `dict`

**uuid**

UUID of the result

**Type** `str`

**extract\_db\_info** (*obj, keys*)

Extract metadata from serialized file

**update\_meta\_info** ()

Extract metadata from myself

`numina.types.structured.load` (*fd*)

” *fd*: file or file-like

File to be opened

`numina.types.structured.open` (*name*)

” *name*: str or file or file-like or `pathlib.Path`

File to be opened

### 3.1.14 `numina.user` — CLI interface

User command line interface of Numina.

### 3.1.15 `numina.util` — Generic utilities

#### `numina.util.flow` — Flow objects

**exception** `numina.util.flow.FlowError`

Error base class for flows.

**class** `numina.util.flow.MixerFlow` (*table*)

**class** `numina.util.flow.ParallelFlow` (*nodeseq*)

A flow where Nodes are executed in parallel.

**class** `numina.util.flow.SerialFlow` (*nodeseq*)

A flow where Nodes are executed sequentially.

#### `numina.util.objimport` — Import objects

Import objects by name

`numina.util.objimport.import_object` (*path*)

Import an object given its fully qualified name.

### 3.1.16 `numina.visualization` — Data visualization

Visualization utilities.

**class** `numina.visualization.ZScaleInterval` (*contrast=0.25*)  
Compute z1 and z2 cuts in a similar way to Iraf.

If the total number of pixels is less than 10, the function simply returns the minimum and the maximum values.

**Parameters** `contrast` (*float, optional*) – The scaling factor (between 0 and 1) for determining the minimum and maximum value. Larger values increase the difference between the minimum and maximum values used for display. Defaults to 0.25.

**:param .. note::** **Deprecated in numina 0.10:** Use `astropy.visualization.ZScaleInterval` instead. It will be removed in numina 1.0

**get\_limits** (*values*)

Return the minimum and maximum value in the interval based on the values provided.

**Parameters** `values` (*~numpy.ndarray*) – The image values.

**Returns** `vmin, vmax` – The minimum and maximum image value in the interval.

**Return type** `float`

### 3.1.17 `numina.user.xdgdirs` — Base Directories

Implementation of some of freedesktop.org Base Directories.

The directories are defined here:

<http://standards.freedesktop.org/basedir-spec/>

We only require `xdg_data_dirs` and `xdg_config_home`

## 3.2 Indices and tables

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- [Glossary](#)

**DFP** Data Factory Pipeline

**DRP** Data Reduction Pipeline

**observing mode** One of the prescribed ways of observing with an instrument

**recipe** A software object that processes the data obtained with a given observing mode of the instrument

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