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# pymcxray Documentation

*Release 0.1.4*

**Hendrix Demers**

**Mar 25, 2021**



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Contents:



# CHAPTER 1

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pymcxray

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Python scripts for using mxray software

- Free software: Apache Software License 2.0
- Documentation: <https://pymcxray.readthedocs.io>.

## 1.1 Features

- TODO

## 1.2 Credits

This package was created with [Cookiecutter](#) and the [audreyr/cookiecutter-pypackage](#) project template.



# CHAPTER 2

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## Installation

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### 2.1 Stable release

---

**Todo:** Add pypi installation for this project.

---

**Warning:** Right now, the *Stable release* installation does not work.

To install pymcxray, run this command in your terminal:

```
$ pip install pymcxray
```

This is the preferred method to install pymcxray, as it will always install the most recent stable release.

If you don't have `pip` installed, this [Python installation guide](#) can guide you through the process.

### 2.2 For developer

Clone the public repository:

```
$ git clone git://github.com/drix00/pymcxray
```

Go in the project folder and install it with pip in developer mode:

```
$ cd pymcxray
$ pip install -e .
```

---

**Note:** The project use Git LFS for the test data file. Follow the information on [Git LFS](#) to get the test data when the repository is pull.

---

## 2.3 From sources

The sources for pymcxray can be downloaded from the Github repo.

You can either clone the public repository:

```
$ git clone git://github.com/drix00/pymcxray
```

Or download the tarball:

```
$ curl -OL https://github.com/drix00/pymcxray/tarball/master
```

Once you have a copy of the source, you can install it with:

```
$ python setup.py install
```

# CHAPTER 3

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## Usage

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To use pymcxray in a project look in the *examples* folder.

To run the example, you need to specify the task you want to do.

```
$python simulation_test_maps.py generate
```

---

**Note:** On windows use the `py` command and not `python` if you have more than one python version installed.

---

Currently these tasks are available:

- generate
- check
- read
- analyze
- scheduled\_read

At first, the script may look complex, but starting with a previous script it is easy to create a new script for your simulation.

Each script has two parts:

- **a class that subclass *mcxray.\_Simulations* were**
  - the simulation parameters are specified
  - results needed extracted from the simulation
  - analysis of the results can be done (this can be also done in another script)
- **a main section were**
  - batch files are generated for the simulation
  - the different task are selected using command line argument.

---

**Note:** These two parts could be separate in two python script (file) without problem. But it seem more logical to kept them together and only have one python script file.

---

## 3.1 simulation\_test\_maps

To explain this example, we are going to start from the main section.

The batch file is define in a run method, which is called by the `__main__`, i.e., only run if the script is called from the command line, but not run if imported.

When the script is ready, these task are run in that order:

1. generate
2. check
3. read
4. analyze

```
$python simulation_test_maps.py generate
```

Start the batch files in parallel. It is easier to just select them all in Windows Explorer and right-click and open them.

```
BatchSimulationTestMapsMM2017_1.bat  
BatchSimulationTestMapsMM2017_2.bat
```

To check the progress of all simulations, i.e., how many simulations are done and todo

```
$python simulation_test_maps.py check
```

When all simulations are done, extract the results and save it in a hdf5 file.

```
$python simulation_test_maps.py read
```

The specified results can be analysed using the script (from the hdf5 file) or analysed using another script. For example, xray-spectrum-modeling project analyse the hdf5 file generate by this script.

An alternative to use the command line argument for these tasks, it to add them directly in the script and uncommenting the task you want to run in the main section of the script.

```
if __name__ == '__main__': #pragma: no cover
    import sys
    ...
    if len(sys.argv) == 1:
        sys.argv.append(mcxray.ANALYZE_TYPE_GENERATE_INPUT_FILE)
        #sys.argv.append(mcxray.ANALYZE_TYPE_CHECK_PROGRESS)
        #sys.argv.append(mcxray.ANALYZE_TYPE_ANALYZE_RESULTS)
        #sys.argv.append(mcxray.ANALYZE_TYPE_ANALYZE_SCHEDULED_READ)
    ...

```

By default the logging level is set at `logging.WARN` and above, but `pymcxray` gives a lot information at the `logging.INFO` level so it is recommended to set the logger level as follow

```
...
logging.getLogger().setLevel(logging.INFO)
...
```

The complete `__main__` is given below

```
if __name__ == '__main__': #pragma: no cover
    import sys
    logging.getLogger().setLevel(logging.INFO)
    logging.info(sys.argv)
    if len(sys.argv) == 1:
        sys.argv.append(mcxray.ANALYZE_TYPE_GENERATE_INPUT_FILE)
        #sys.argv.append(mcxray.ANALYZE_TYPE_CHECK_PROGRESS)
        #sys.argv.append(mcxray.ANALYZE_TYPE_ANALYZE_RESULTS)
        #sys.argv.append(mcxray.ANALYZE_TYPE_ANALYZE_SCHEDULED_READ)
    run()
```

The `run` method have three components

- a configuration file, see [Configuration file](#)
- a batch file object, see [Batch file](#)
- one or more simulation subclass, see [Simulation subclass](#)

Here is a example of a complete `run` method

```
def run():
    # import the batch file class.
    from pymcxray.BatchFileConsole import BatchFileConsole

    # Find the configuration file path
    configuration_file_path = get_current_module_path(__file__, "MCXRay_latest.cfg")
    program_name = get_mcxray_program_name(configuration_file_path)

    # Create the batch file object.
    batch_file = BatchFileConsole("BatchSimulationTestMapsMM2017", program_name,
                                   numberFiles=10)

    # Create the simulation object and add the batch file object to it.
    analyze = SimulationTestMapsMM2017(relativePath=r"mcxray/SimulationTestMapsMM2017",
                                         configurationFilepath=configuration_file_path)
    analyze.run(batch_file)
```

## 3.2 Configuration file

The configuration file is a ini style configuration file for using pymcxray. It define the paths needed to generate and run the simulations.

```
[Paths]
mcxrayProgramName=console_mcxray_x64.exe
resultsMcGillPath=D:\Dropbox\hdemers\professional\results\simulations
mcxrayArchivePath=D:\Dropbox\hdemers\professional\softwareRelease\mcxray
mcxrayArchiveName=2016-04-11_11h41m28s_MCXRay_v1.6.6.0.zip
```

See the documentation of these functions for more detail on each option

- `pymcxray.get_mcxtay_program_name()`
- `pymcxray.get_results_mcgill_path()`
- `pymcxray.get_mcxtay_program_path()`
- `pymcxray.get_mcxtay_archive_path()`

### 3.3 Batch file

The batch file is responsible to create the simulation structure with a copy of mcxtay program. Batch files are generated to easily run the simulations. One important parameter to set is the `numberFiles`, this is the number of batch files generated and that can be run in parallel. For maximum efficiency it should be set as the number of logical processors minus 1 or 2. For example, on a computer with 12 logical processors, the `numberFiles` should be set at 10.

See `pymcxray.BatchFileConsole.BatchFileConsole` documentation for more information about the other parameters.

### 3.4 Simulation subclass

The simulation subclass allows to generate a lot of simulations by varying simulations parameters.

The main features are

- **generate input files**
  - regenerate input files of simulations not done
- check the progress of the simulations: done and todo
- **extract results from the completed simulation**
  - save the results in a hdf5 file for easier analysis
- optionally do the analysis of the results

To do that the user need to subclass `pymcxray.mcxtay._Simulations` and overwrite these method

- `pymcxray.mcxtay._Simulations._initData()` (required)
- `pymcxray.mcxtay._Simulations.getAnalysisName()` (required)
- `pymcxray.mcxtay._Simulations.createSpecimen()` (required)
- `pymcxray.mcxtay._Simulations.read_one_results_hdf5()` (optional)
- `pymcxray.mcxtay._Simulations.analyze_results_hdf5()` (optional)

**Warning:** If any of the required method is modified, the simulation have to be redone completely. It is recommended to just delete the root path for the analysis and generate the input files and do the simulations. For this example, delete `SimulationTestMapsMM2017` folder.

**Warning:** If `pymcxray.mcxtay._Simulations.read_one_results_hdf5()` is modified. In some case, the hdf5 need to be deleted. Furthermore, if the results were deleted: `delete_result_files` is `True`, the simulation have to be redone.

Below are given example for each method, for more detail see the method documentation.

### 3.4.1 Init data

This method is used to specify the options for the analysis and also the parameters used in the simulations.

The most important options for the analysis are:

- *use\_hdf5* to use the recommended hdf5 method. If it is *False* the older serial method will be used.
- *delete\_result\_files* if it is *True*, the result file are deleted after added in the hdf5 file. Very useful when creating a lot files like for a map.
- *createBackup* if *True* create a backup of the hdf5 file before adding more results to it. Usefull to not loss data in case of error or crash, but you should delete backup file manually as they can take a lot of space.

**Warning:** If *delete\_result\_files* is *True*, all results are deleted for a simulation and only the results specified in `pymcxray.mcxray._Simulations.read_one_results_hdf5()` are kept. If `pymcxray.mcxray._Simulations.read_one_results_hdf5()` is modified to extract more results, the simulation have to be simulate again.

This is the recommended values for the options, only change them when you are sure everything is working OK.

```
def __initData(self):
    self.use_hdf5 = True
    self.delete_result_files = False
    self.createBackup = True
```

The simulation parameters are specified in this method.

---

**Note:** If not specified, the script use MCXRay default parameters. Start MCXRay program to see the default value of each parameters.

---

To change simulation parameters, create a `pymcxray.SimulationsParameters.SimulationsParameters` object

```
self._simulationsParameters = SimulationsParameters()
```

Two kinds of parameter can be added:

- varied specified with a list of values
- fixed specified with a single value

The script will automatically generate a simulation for all combination of the varied parameters.

**Warning:** Adding a lot of varied parameters with a lot of values can generate a lot of simulations. Above 1000 simulations, all tasks of the script will be slow because of the generation or reading of a lot of files. It is recommended to start with 2 or 3 varied parameters and short list of values and test all tasks of the script. When the tests are OK and results make sense, you can increase the list of values. To add more varied parameters, it is recommended to create a new script with again only 2 or 3 varied parameters.

The parameters that can be added are defined as keyword starting with `PARAMETER_` in the module `pymcxray.SimulationsParameters`. If you don't find the parameter you want request an "enhancement" at <https://github.com/drix00/pymcxray/issues>.

Here is an example how-to add simulation parameters

```
from pymcxray.SimulationsParameters import SimulationsParameters, PARAMETER INCIDENT_
    ENERGY_keV, PARAMETER_NUMBER_ELECTRONS, \
PARAMETER_BEAM_POSITION_nm, PARAMETER_NUMBER_XRAYS
...
class SimulationTestMapsMM2017(mcxray._Simulations):
    def __initData(self):
        ...

        # Local variables for value and list if values.
        energy_keV = 30.0
        number_electrons = 10000

        #number_xrays_list = [10, 20, 30, 50, 60, 100, 200, 500, 1000]
        number_xrays_list = [10]
        xs_nm = np.linspace(-5.0e3, 5.0e3, 3)
        probePositions_nm = [tuple(position_nm) for position_nm in np.transpose([np.
            tile(xs_nm, len(xs_nm)), np.repeat(xs_nm, len(xs_nm))]).tolist()]

        # Simulation parameters
        self._simulationsParameters = SimulationsParameters()

        self._simulationsParameters.addVaried(PARAMETER_NUMBER_XRAYS, number_xrays_
list)
        self._simulationsParameters.addVaried(PARAMETER_BEAM_POSITION_nm,_
probePositions_nm)

        self._simulationsParameters.addFixed(PARAMETER INCIDENT ENERGY_keV, energy_-
keV)
        self._simulationsParameters.addFixed(PARAMETER_NUMBER_ELECTRONS, number_-
electrons)
```

### 3.4.2 Analysis name

This method specify the name of the analysis or experiment for which the simulation are done. Normally similar to the name of the class and mostly used as basename for the input files and result files. In case of two scripts writing the same path, it allows to differentiate them, but it is not recommended to run two scripts in the same folde.

```
class SimulationTestMapsMM2017(mcxray._Simulations):
    ...
    def getAnalysisName(self):
        return "SimulationTestMapsMM2017"
    ...
```

### 3.4.3 Create specimen

This method is used to create the specimen for each simulation. The argument of the method contains the option of the specific simulation and can be used to create the specimen. The `pymcxray.Simulation.Simulation` module contains predefined specimen which can be use in this method.

Here an example how-to use the *parameters* argument and the predefined specimen

```
def createSpecimen(self, parameters):
    weightFractions = parameters[PARAMETER_WEIGHT_FRACTIONS]

    elements = [(self.atomicNumberA, weightFractions[0]),
                (self.atomicNumberB, weightFractions[1])]
    specimen = Simulation.createAlloyBulkSample(elements)
    return specimen
```

A more complex example, where each region are specified is given below

```
def createSpecimen(self, parameters):
    # Create the specimen with a name and number of regions.
    specimen = Specimen.Specimen()
    specimen.name = "Maps01"
    specimen.numberRegions = 10

    # Region 0
    region = Region.Region()
    region.numberElements = 0
    region.regionType = RegionType.REGION_TYPE_BOX
    parameters = [-10000000000.0, 10000000000.0, -10000000000.0, 10000000000.0, 0.0,
    ↳20000000000.0]
    region.regionDimensions = RegionDimensions.RegionDimensionsBox(parameters)
    specimen.regions.append(region)

    # Region 1
    region = Region.Region()
    region.numberElements = 2
    region.elements = [Element.Element(27, massFraction=0.01), Element.Element(26,
    ↳massFraction=0.99)]
    region.regionType = RegionType.REGION_TYPE_BOX
    parameters = [-7.5e4, -2.5e4, -7.5e4, -2.5e4, 0.0, 0.2e4]
    region.regionDimensions = RegionDimensions.RegionDimensionsBox(parameters)
    specimen.regions.append(region)
    ...
```

**Warning:** Creating a specimen in MCXRay is complicate as sometime you need to create a empty region 0. Look at the predefined specimen in `pymcxray.Simulation.Simulation` for help. Drawing the trajectory with the `FileFormat.Results.ElectronTrajectoriesResults` will help debug the specimen. You can also request a “help wanted” at <https://github.com/drix00/pymcxray/issues>.

### 3.4.4 Read one simulation results

This method extract results from one complete simulation and added them in a hdf5 group for this simulation. The result that can be extracted are in the package `pymcxray.FileFormat.Results` and only the class implementing `write_hdf5()` can be extracted. If the desired results does not implement the `write_hdf5()` method, request an “enhancement” at <https://github.com/drix00/pymcxray/issues>.

---

**Note:** The format of the hdf5 file is not well documented. Check the implementation of the `write_hdf5()` method and request an “enhancement” at <https://github.com/drix00/pymcxray/issues> for the documentation.

---

**Note:** The program HDFView is useful to look at the hdf5 file. See <https://support.hdfgroup.org/products/java/hdfview/>.

---

Here is an example how-to extract the electron results (BSE, TE, ...)

```
def read_one_results_hdf5(self, simulation, hdf5_group):
    electronResults = ElectronResults.ElectronResults()
    electronResults.path = self.getSimulationsPath()
    electronResults.basename = simulation.resultsBasename
    electronResults.read()
    electronResults.write_hdf5(hdf5_group)
```

So far this class are implemented with hdf5 support

- `pymcxray.FileFormat.Results.ElectronResults.ElectronResults`
- `pymcxray.FileFormat.Results.PhirhozEmittedCharacteristic.PhirhozEmittedCharacteristic`
- `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristic.PhirhozGeneratedCharacteristic`
- `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristicThinFilm.PhirhozGeneratedCharacteristicThinFilm`
- `pymcxray.FileFormat.Results.XrayIntensities.XrayIntensities`
- `pymcxray.FileFormat.Results.XraySpectraRegionsEmitted.XraySpectraRegionsEmitted`
- `pymcxray.FileFormat.Results.XraySpectraSpecimenEmittedDetected.XraySpectraSpecimenEmittedDetected`

### 3.4.5 Analyze all simulations

This method is only needed for the task analyze.

Often the method start by calling `mcxray._Simulations.readResults()` to read all new results and add them in the hdf5 file.

The example below shows how-to open the hdf5 file in memory for somewhat fast analysis simulation. The file is read only at the beginning and stored in memory.

```
def analyze_results_hdf5(self): #pragma: no cover
    self.readResults()

    file_path = self.get_hdf5_file_path()
    with h5py.File(file_path, 'r', driver='core') as hdf5_file:
        hdf5_group = self.get_hdf5_group(hdf5_file)
        logging.info(hdf5_group.name)
```

---

**Todo:** Document `pymcxray.mcxray`

---

---

**Todo:** Document `pymcxray.SimulationsParameters`

---

---

**Todo:** Document `pymcxray.Simulation`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.ElectronResults.ElectronResults`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.PhirhozEmittedCharacteristic.PhirhozEmittedCharacteristic`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristic.PhirhozGeneratedCharacteristic`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristicThinFilm.PhirhozGeneratedCharacteristicThinFilm`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.XrayIntensities.XrayIntensities`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.XraySpectraRegionsEmitted.XraySpectraRegionsEmitted`

---

---

**Todo:** Document `pymcxray.FileFormat.Results.XraySpectraSpecimenEmittedDetected.XraySpectraSpecimenEmittedDetected`

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# CHAPTER 4

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## Contributing

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Contributions are welcome, and they are greatly appreciated! Every little bit helps, and credit will always be given. You can contribute in many ways:

### 4.1 Types of Contributions

#### 4.1.1 Report Bugs

Report bugs at <https://github.com/drix00/pymcxray/issues>.

If you are reporting a bug, please include:

- Your operating system name and version.
- Any details about your local setup that might be helpful in troubleshooting.
- Detailed steps to reproduce the bug.

#### 4.1.2 Fix Bugs

Look through the GitHub issues for bugs. Anything tagged with “bug” and “help wanted” is open to whoever wants to implement it.

#### 4.1.3 Implement Features

Look through the GitHub issues for features. Anything tagged with “enhancement” and “help wanted” is open to whoever wants to implement it.

#### 4.1.4 Write Documentation

pymcxray could always use more documentation, whether as part of the official pymcxray docs, in docstrings, or even on the web in blog posts, articles, and such.

#### 4.1.5 Submit Feedback

The best way to send feedback is to file an issue at <https://github.com/drix00/pymcxray/issues>.

If you are proposing a feature:

- Explain in detail how it would work.
- Keep the scope as narrow as possible, to make it easier to implement.
- Remember that this is a volunteer-driven project, and that contributions are welcome :)

## 4.2 Get Started!

Ready to contribute? Here's how to set up *pymcxray* for local development.

1. Fork the *pymcxray* repo on GitHub.

2. Clone your fork locally:

```
$ git clone git@github.com:your_name_here/pymcxray.git
```

3. Install your local copy into a virtualenv. Assuming you have `virtualenvwrapper` installed, this is how you set up your fork for local development:

```
$ mkvirtualenv pymcxray
$ cd pymcxray/
$ python setup.py develop
```

4. Create a branch for local development:

```
$ git checkout -b name-of-your-bugfix-or-feature
```

Now you can make your changes locally.

5. When you're done making changes, check that your changes pass flake8 and the tests, including testing other Python versions with tox:

```
$ flake8 pymcxray tests
$ python setup.py test or py.test
$ tox
```

To get flake8 and tox, just pip install them into your virtualenv.

6. Commit your changes and push your branch to GitHub:

```
$ git add .
$ git commit -m "Your detailed description of your changes."
$ git push origin name-of-your-bugfix-or-feature
```

7. Submit a pull request through the GitHub website.

## 4.3 Pull Request Guidelines

Before you submit a pull request, check that it meets these guidelines:

1. The pull request should include tests.
2. If the pull request adds functionality, the docs should be updated. Put your new functionality into a function with a docstring, and add the feature to the list in README.rst.
3. The pull request should work for Python 2.6, 2.7, 3.3, 3.4 and 3.5, and for PyPy. Check [https://travis-ci.org/drix00/pymcxray/pull\\_requests](https://travis-ci.org/drix00/pymcxray/pull_requests) and make sure that the tests pass for all supported Python versions.

## 4.4 Tips

To run a subset of tests:

```
$ python -m unittest tests.test_pymcxray
```



# CHAPTER 5

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## Credits

---

### 5.1 Development Lead

- Hendrix Demers <[hendrix.demers@mail.mcgill.ca](mailto:hendrix.demers@mail.mcgill.ca)>

### 5.2 Contributors

None yet. Why not be the first?



# CHAPTER 6

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## History

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### 6.1 0.1.4 (2019-11-29)

- Add calculation of weight fractions for trace elements.

### 6.2 0.1.2 (2017-06-26)

- Add create multi-layers specimen.
- Add number of layers X, Y, and Z as parameters.

### 6.3 0.1.0 (2017-03-07)

- First release on PyPI.



# CHAPTER 7

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## ToDo

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**Todo:** Add pypi installation for this project.

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/installation.rst, line 13.)

---

**Todo:** Document `pymcxray.mcxray`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 436.)

---

**Todo:** Document `pymcxray.SimulationsParameters`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 437.)

---

**Todo:** Document `pymcxray.Simulation`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 438.)

---

**Todo:** Document `pymcxray.FileFormat.Results.ElectronResults.ElectronResults`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 439.)

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**Todo:** Document `pymcxray.FileFormat.Results.PhirhozEmittedCharacteristic`.  
`PhirhozEmittedCharacteristic`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 440.)

---

**Todo:** Document `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristic`.  
`PhirhozGeneratedCharacteristic`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 441.)

---

**Todo:** Document `pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristicThinFilm`.  
`PhirhozGeneratedCharacteristicThinFilm`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 442.)

---

**Todo:** Document `pymcxray.FileFormat.Results.XrayIntensities`.  
`XrayIntensities`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 443.)

---

**Todo:** Document `pymcxray.FileFormat.Results.XraySpectraRegionsEmitted`.  
`XraySpectraRegionsEmitted`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 444.)

---

**Todo:** Document `pymcxray.FileFormat.Results.XraySpectraSpecimenEmittedDetected`.  
`XraySpectraSpecimenEmittedDetected`

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/docs/usage.rst, line 445.)

---

**Todo:** Add units.

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/pymcxray/ElementProperties of pymcxray.ElementProperties.g\_FermiEnergy, line 6.)

---

**Todo:** Add units.

---

(The original entry is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/pymcxray/ElementProperties of pymcxray.ElementProperties.g\_kFermi, line 6.)

---

**Todo:** Add units.

---

(The [original entry](#) is located in /home/docs/checkouts/readthedocs.org/user\_builds/pymcxray/checkouts/latest/pymcxray/ElementProperties.g\_phasmonEnergy, line 6.)



# CHAPTER 8

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pymcxray

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## 8.1 pymcxray package

### 8.1.1 Subpackages

**pymcxray.FileFormat package**

**Subpackages**

**pymcxray.FileFormat.Results package**

**Subpackages**

**pymcxray.FileFormat.Results.exported package**

**Submodules**

**pymcxray.FileFormat.Results.exported.DataMap module**

Read data map exported by MCXRay.

```
class pymcxray.FileFormat.Results.exported.DataMap(filepath)
    Bases: object

    imageName
    pixels
    read()
    saveImage(imageFilepath=None)
    showImage()
```

**size**

```
pymcxray.FileFormat.Results.exported.DataMap.run()
```

## pymcxray.FileFormat.Results.exported.XrayIntensityXY module

Read x-ray intensity distribution in XY exported from mcxray GUI.

```
class pymcxray.FileFormat.Results.exported.XrayIntensityXY.XrayIntensityXY
```

Bases: object

```
readData(filepath)
```

## pymcxray.FileFormat.Results.exported.test\_DataMap module

Tests for the module *DataMap*.

```
class pymcxray.FileFormat.Results.exported.test_DataMap.TestDataMap(methodName='runTest')
```

Bases: unittest.case.TestCase

TestCase class for the module *DataMap*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
```

Setup method.

```
tearDown()
```

Teardown method.

```
testSkeleton()
```

First test to check if the testcase is working with the testing framework.

```
test_init()
```

Tests for method *init*.

```
test_read()
```

Tests for method *read*.

## pymcxray.FileFormat.Results.exported.test\_XrayIntensityXY module

Tests for the module *XrayIntensityXY*.

```
class pymcxray.FileFormat.Results.exported.test_XrayIntensityXY.TestXrayIntensityXY(methodName='runTest')
```

Bases: unittest.case.TestCase

TestCase class for the module *XrayIntensityXY*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
```

Setup method.

```
tearDown()
```

Teardown method.

```
testOpenFile()
Test if the test data file can be open.

testSkeleton()
First test to check if the testcase is working with the testing framework.
```

## pymcxray.FileFormat.Results.exported.tests module

### Module contents

#### Submodules

## pymcxray.FileFormat.Results.BaseResults module

BaseResults

```
class pymcxray.FileFormat.Results.BaseResults.BaseResults(path='',           base-
                                                       name='MCXRay')
Bases: object

basename
filepath
path
```

## pymcxray.FileFormat.Results.BeamParameters module

MCXRay beam parameters from results file.

```
class pymcxray.FileFormat.Results.BeamParameters.BeamParameters
Bases: object

acquisitionTime_s
current_A
diameter90_A
gaussianMean
gaussianSigma
incidentEnergy_keV
readFromLines(lines)
tiltAngle_deg
```

## pymcxray.FileFormat.Results.DetectorParameters module

MCXRay detector parameters from results file.

```
class pymcxray.FileFormat.Results.DetectorParameters.DetectorParameters
Bases: object

angleBetweenDetectorSpecimenNormal_deg
angleBetweenDetectorXAxis_deg
```

```
beamDetectorDistance_cm
crystalDensity_g_cm3
crystalName
crystalRadius_cm
crystalThickness_cm
deadLayerThickness_A
diffusionLength_A
noiseEdsDetector_eV
readFromLines(lines)
solidAngle_deg
surfaceQualityFactor
takeoffAngleEffective_deg
takeoffAngleNormalIncidence_deg
thicknessAir_um
thicknessAlWindow_um
thicknessBeWindow_um
thicknessH2O_um
thicknessMoxtek_um
thicknessOil_um
thicknessTiWindow_um
```

## pymcxray.FileFormat.Results.Dump module

MCXRay dump results file.

```
class pymcxray.FileFormat.Results.Dump.Dump
    Bases: object
    read(filepath)
```

## pymcxray.FileFormat.Results.ElectronExistResults module

Read electron exit results simulated with MCXRay.

```
class pymcxray.FileFormat.Results.ElectronExistResults.ElectronDetector
    Bases: object
    azimuthalAngleRange_deg
    detectElectrons(data)
    energyRange_keV
    maximumAzimuthalAngle_deg
    maximumEnergy_keV
```

```
maximumPolarAngle_deg
minimumAzimuthalAngle_deg
minimumEnergy_keV
minimumPolarAngle_deg
polarAngleRange_deg

class pymcxray.FileFormat.Results.ElectronExistResults(*args,
                                                       **kargs)
    Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults
    data
        getEnergyDistribution(numberBins=10)
        numberData
        read()
```

## pymcxray.FileFormat.Results.ElectronParameters module

MCXRay electron parameters results file.

```
class pymcxray.FileFormat.Results.ElectronParameters.ElectronParameters
    Bases: object
    backscatteredRatio
    eRatio
    internalRatio
    meanAzimuthalAngleCollision_deg
    meanDistanceBetweenCollisions_A
    meanNumberCollisionPerElectrons
    meanPolarAngleCollision_deg
    numberSimulatedElectrons
    readFromLines(lines)
    skirtRatio
    throughRatio
```

## pymcxray.FileFormat.Results.ElectronResults module

Read ElectronResults MCXRay results file.

```
class pymcxray.FileFormat.Results.ElectronResults.ElectronResults
    Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults
    fieldNames
        fractionBackscatteredElectrons
        fractionInternalElectrons
        fractionSkirtedElectrons
```

```
fractionTransmittedElectrons
numberBackscatteredElectrons
numberElectronCollisions
numberInternalElectrons
numberSimulatedElectrons
numberSkirtedElectrons
numberTransmittedElectrons
read()
write_hdf5(hdf5_group)
```

## pymcxray.FileFormat.Results.ElectronTrajectoriesResults module

Read MCXray electron trajectories results file.

```
class pymcxray.FileFormat.Results.ElectronTrajectoriesResults.Collision
Bases: object

    collisionType
    correctedX_A
    correctedY_A
    correctedZ_A
    energy_keV
    indexRegion
    x_A
    y_A
    z_A

class pymcxray.FileFormat.Results.ElectronTrajectoriesResults.ElectronTrajectoriesResults()
Bases: object

    drawXY(title='', corrected=False, x_limit=None, y_limit=None)
    drawXZ(title='', corrected=False, theta_deg=0.0, colorType='colorTrajectoryType', trajectoryIndexes=None, x_limit=None, y_limit=None)
    drawYZ(title='', corrected=False, x_limit=None, y_limit=None)
    getElectronGunPositions_nm()
    read(filepath)
    write_hdf5(hdf5_group)

class pymcxray.FileFormat.Results.ElectronTrajectoriesResults.Trajectory
Bases: object

    addCollision(collision)
    collisions
    index
```

**trajectoryType**

```
pymcxray.FileFormat.Results.ElectronTrajectoriesResults.run()  
pymcxray.FileFormat.Results.ElectronTrajectoriesResults.runAuCThinFilm()  
pymcxray.FileFormat.Results.ElectronTrajectoriesResults.runFogging()
```

**pymcxray.FileFormat.Results.ElementParameters module**

MCXRay element parameters result file.

```
class pymcxray.FileFormat.Results.ElementParameters.ElementParameters  
Bases: object
```

**pymcxray.FileFormat.Results.Intersections module**

MCXRay intersections results file.

```
class pymcxray.FileFormat.Results.Intersections.Intersections  
Bases: object  
extractFromLines (lines)
```

**pymcxray.FileFormat.Results.MicroscopeParameters module**

MCXRay microscope parameter in results file.

```
class pymcxray.FileFormat.Results.MicroscopeParameters.MicroscopeParameters  
Bases: object  
beamParameters  
detectorParameters  
readFromLines (lines)
```

**pymcxray.FileFormat.Results.ModelParameters module**

MCXRay model parameters from results file.

```
class pymcxray.FileFormat.Results.ModelParameters.ModelParameters  
Bases: object  
atomCollisionModel  
atomCollisionScreeningModel  
atomCrossSectionModel  
atomCrossSectionScreeningModel  
atomElectronRangeModel  
atomEnergyLossModel  
atomMeanIonizationPotentialModel  
atomScreeningModel
```

```
bremsstrahlungCrossSectionModel
characteristicCrossSectionModel
readFromLines (lines)
regionEnergyLossModel
```

## pymcxray.FileFormat.Results.Phirhoz module

MCXRay phirhoz result file.

```
class pymcxray.FileFormat.Results.Phirhoz.Phirhoz (symbol, shell)
Bases: object

depths_A
intensity
readFromLines (lines)
shell
symbol
values
```

## pymcxray.FileFormat.Results.PhirhozElement module

MCXRay result file phirhoz element.

```
class pymcxray.FileFormat.Results.PhirhozElement.PhirhozElement
Bases: object

isIonizationShell_K
isIonizationShell_L
isIonizationShell_M
readFromLine (line)
symbol
weightFraction
```

## pymcxray.FileFormat.Results.PhirhozEmittedCharacteristic module

Read PhirhozEmittedCharacteristic file from MCXRay.

```
class pymcxray.FileFormat.Results.PhirhozEmittedCharacteristic.PhirhozEmittedCharacteristic
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

depth_A
depth_nm
fieldNames
phirhozs
read (regionID=0)
```

```
write_hdf5 (hdf5_group)
```

## pymcxray.FileFormat.Results.PhirhozEmittedCharacteristicThinFilm module

Read mcxray phirhoz thin film emitted results.

```
class pymcxray.FileFormat.Results.PhirhozEmittedCharacteristicThinFilm.PhirhozEmittedCharacteristicThinFilm
    Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

    fieldNames
    getIntensity (regionID, atomicSymbol, xrayLine)
    intensities
    numberRegions
    read()
```

## pymcxray.FileFormat.Results.PhirhozGenerated module

MCXRay phirhoz generated result file.

```
class pymcxray.FileFormat.Results.PhirhozGenerated.PhirhozGenerated
    Bases: object

    electronParameters
    getCharacteristicPhiRhoZ (regionID)
    microscopeParameters
    modelParameters
    numberRegions
    read (filepath)
    readElectron (lines)
    readMicroscope (lines)
    readRegions (lines)
    readSimulationParameters (lines)
    readSolidSimulationModels (lines)
    simulationParameters
```

## pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristic module

Read PhirhozGeneratedCharacteristic file from MCXRay program.

```
class pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristic.PhirhozGeneratedCharacteristic
    Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

    depth_A
    depth_nm
    fieldNames
```

```
phirhozs
read(regionID=0)
write_hdf5(hdf5_group)
```

## pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristicThinFilm module

Read mcxray phirhoz thin film generated results.

```
class pymcxray.FileFormat.Results.PhirhozGeneratedCharacteristicThinFilm.PhirhozGeneratedCharacteristicThinFilm
    Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

    fieldNames
    getIntensity(regionID, atomicSymbol, xrayLine)
    get_subshells()
    get_symbols()
    intensities
    numberRegions
    read()
    write_hdf5(hdf5_group)
```

## pymcxray.FileFormat.Results.PhirhozRegion module

MCXRay phirhoz results file for a region.

```
class pymcxray.FileFormat.Results.PhirhozRegion.PhirhozRegion(numberEnergyWindows,
                                                               numberLayersZ)
    Bases: object

    readBackgroundPhirhozs(lines)
    readCharacteristicPhirhozs(lines)
    readElements(lines)
    readFromLines(lines)
    readPhirhozDistributions(lines)
    readRegionVolume(lines)
    regionID
```

## pymcxray.FileFormat.Results.RegionParameters module

MCXRay region parameters result file.

```
class pymcxray.FileFormat.Results.RegionParameters.RegionParameters
    Bases: object

    elements
    layerThickness_A
    numberElements
```

```
regionID
```

## pymcxray.FileFormat.Results.RegionVolume module

MCXRay result file region volume.

```
class pymcxray.FileFormat.Results.RegionVolume.RegionVolume
Bases: object
    readFromLines (lines)
    regionID
```

## pymcxray.FileFormat.Results.SimulationParameters module

MCXRay simulation parameters results file.

```
class pymcxray.FileFormat.Results.SimulationParameters.SimulationParameters
Bases: object
    edsMaximumEnergy_keV
    generalizedWalk
    interpolationType
    maximumLiveTime_s
    numberChannels
    numberElectrons
    numberEnergyWindows
    numberLayersX
    numberLayersY
    numberLayersZ
    numberPhotons
    readFromLines (lines)
    useLiveTime_s
```

## pymcxray.FileFormat.Results.Spectra module

MCXRay spectra result file.

```
class pymcxray.FileFormat.Results.Spectra.Spectra
Bases: object
    getElementSpectra()
    getElementSpectrum (regionID, elementName)
    getRegionParameters (regionID)
    getRegionSpectrum (regionID)
    getSpecimenSpectrum()
```

```
numberRegions
read(filepath)
readRegion(lines)
readRegions(lines)
readSpecimen(lines)
```

## pymcxray.FileFormat.Results.SpectraEDS module

Read MCXRay spectra EDS results file.

```
class pymcxray.FileFormat.Results.SpectraEDS.SpectraEDS
    Bases: object

    characteristicProbability
    elements
    numberCharacteristicPeaks
    numberElements
    numberSimulatedPhotons
    readFileObject(inputFile)
    readfilepath(filepath)
    readPartialSpectraReferenceSection(lines)
    readRegionSpectraSection(lines)
    readTestInputSection(lines)
    regionID

pymcxray.FileFormat.Results.SpectraEDS.runExample()
```

## pymcxray.FileFormat.Results.Spectrum module

MCXRay spectrum result file.

```
class pymcxray.FileFormat.Results.Spectrum.Spectrum
    Bases: object

    backgroundIntensities
    characteristicIntensities
    energies_keV
    intensities
```

## pymcxray.FileFormat.Results.SpectrumEDS module

Read MCXRay EDS spectrm from results file.

```
class pymcxray.FileFormat.Results.SpectrumEDS.SpectrumEDS(lines=None)
    Bases: object
```

```
channels
countsList
enegies_keV
readLines (lines)
```

## pymcxray.FileFormat.Results.Tags module

MCXRay tags used in output files.

```
exception pymcxray.FileFormat.Results.Tags.TagNotFoundError
Bases: ValueError

pymcxray.FileFormat.Results.Tags.findallTag (tag, lines, contains=None)
pymcxray.FileFormat.Results.Tags.findTag (tag, lines)
```

## pymcxray.FileFormat.Results.XrayIntensities module

Xray intensities result file from MCXRay.

```
class pymcxray.FileFormat.Results.XrayIntensities.XrayIntensities
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

fieldNames
getAtomicNumberLineEnergySets ()
getDetectedIntensity (atomicNumber, xrayLine)
getIntensity (data_type, region_id, atomic_number, xray_line)
getIntensityEmitted (atomicNumber, xrayLine, total=True)
getIntensityEmittedDetected (atomicNumber, xrayLine, total=True)
getIntensityGenerated (atomicNumber, xraySubshell, total=False)
getIntensityGeneratedDetected (atomicNumber, xraySubshell)
getResult_types ()
getXray_lines ()
intensities
numberIntensities
read ()
write_hdf5 (hdf5_group)
```

## pymcxray.FileFormat.Results.XraySimulatedSpectraRegion module

description

```
class pymcxray.FileFormat.Results.XraySimulatedSpectraRegion.XraySimulatedSpectraRegion
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

channelNumbers
```

```
detectedIntensities
eNetPeak
energiesReference_keV
energies_keV
fieldNames
peakToBackground
peakToBackgroundAverage
read(regionID=0)
simulatedIntensities
```

## pymcxray.FileFormat.Results.XraySimulatedSpectraSpecimen module

description

```
class pymcxray.FileFormat.Results.XraySimulatedSpectraSpecimen.XraySimulatedSpectraSpecimen
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults
energies_keV
fieldNames
read()
totals
```

## pymcxray.FileFormat.Results.XraySpectraAtomEmittedDetectedLines module

Read MCXRay XraySpectraAtomEmittedDetectedLines file.

```
class pymcxray.FileFormat.Results.XraySpectraAtomEmittedDetectedLines.XraySpectraAtomEmitted
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults
characteristics
energies_keV
fieldNames
read(regionID)
```

## pymcxray.FileFormat.Results.XraySpectraRegionEmitted module

Read XraySpectraRegionEmitted mcxray result file.

```
class pymcxray.FileFormat.Results.XraySpectraRegionEmitted.XraySpectraRegionEmitted
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults
bremsstrahlungValue_1_ekeVsr(energy_keV)
bremsstrahlung_1_ekeVsr
characteristicValue_1_ekeVsr(energy_keV)
characteristic_1_ekeVsr
```

```
energies_keV
fieldnames
read(regionID=0)
totalValue_1_ekeVsr(energy_keV)
total_1_ekeVsr

pymcxray.FileFormat.Results.XraySpectraRegionEmitted.run()
```

## pymcxray.FileFormat.Results.XraySpectraRegionsEmitted module

Read all XraySpectraRegionsEmitted mcxray result files for all regions.

```
class pymcxray.FileFormat.Results.XraySpectraRegionsEmitted.XraySpectraRegionsEmitted
Bases: pymcxray.FileFormat.Results.XraySpectraRegionEmitted.
XraySpectraRegionEmitted

read()
write_hdf5(hdf5_group)

pymcxray.FileFormat.Results.XraySpectraRegionsEmitted.run()
```

## pymcxray.FileFormat.Results.XraySpectraSpecimen module

Read XraySpectraSpecimen MCXRay results file.

```
class pymcxray.FileFormat.Results.XraySpectraSpecimen.XraySpectraSpecimen
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

backgrounds
characteristics
energies_keV
fieldNames
read()
totals
write_hdf5(hdf5_group)
```

## pymcxray.FileFormat.Results.XraySpectraSpecimenEmittedDetected module

Read XraySpectraSpecimenEmittedDetected MCXRay results file.

```
class pymcxray.FileFormat.Results.XraySpectraSpecimenEmittedDetected.XraySpectraSpecimenEmittedDetected
Bases: pymcxray.FileFormat.Results.BaseResults.BaseResults

backgrounds
characteristics
energies_keV
fieldNames
```

```
read()
totals
write_hdf5(hdf5_group)
```

## pymcxray.FileFormat.Results.test\_BaseResults module

Tests for the module *BaseResults*.

```
class pymcxray.FileFormat.Results.test_BaseResults.TestBaseResults(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *BaseResults*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_basename()
    Tests for method basename.

test_init()
    Tests for method init.

test_path()
    Tests for method path.
```

## pymcxray.FileFormat.Results.test\_BeamParameters module

Tests for module *BeamParameters*.

```
class pymcxray.FileFormat.Results.test_BeamParameters.TestBeamParameters(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *BeamParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_readFromLines()
    Tests for method readFromLines.
```

## pymcxray.FileFormat.Results.test\_DetectorParameters module

Tests for module *DetectorParameters*.

**class** pymcxray.FileFormat.Results.test\_DetectorParameters.**TestDetectorParameters** (*methodName*=*runTest*)  
Bases: unittest.case.TestCase

TestCase class for the module *DetectorParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_readFromLines()**

Tests for method *readFromLines*.

## pymcxray.FileFormat.Results.test\_Dump module

Tests for module *Dump*

**class** pymcxray.FileFormat.Results.test\_Dump.**TestDump** (*methodName*=*'runTest'*)  
Bases: unittest.case.TestCase

TestCase class for the module *Dump*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_ElectronExistResults module

### pymcxray.FileFormat.Results.test\_ElectronParameters module

Tests for module *ElectronParameters*.

**class** pymcxray.FileFormat.Results.test\_ElectronParameters.**TestElectronParameters** (*methodName*=*'runTest'*)  
Bases: unittest.case.TestCase

TestCase class for the module *ElectronParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_readFromLines()**

Tests for method *readFromLines*.

```
pymcxray.FileFormat.Results.test_ElectronParameters.getLinesAndReference()
```

## pymcxray.FileFormat.Results.test\_ElectronResults module

Tests for the module *XraySpectraSpecimen*.

```
class pymcxray.FileFormat.Results.test_ElectronResults.TestElectronResults(methodName='runTest')
    Bases: unittest.case.TestCase
```

TestCase class for the module *ElectronResults*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_ElementParameters module

Tests for the module *ElementParameters*.

```
class pymcxray.FileFormat.Results.test_ElementParameters.TestElementParameters(methodName='runTest')
    Bases: unittest.case.TestCase
```

TestCase class for the module *ElementParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

## pymcxray.FileFormat.Results.test\_Intersections module

Tests for module *Intersections*.

**class** pymcxray.FileFormat.Results.test\_Intersections.**TestIntersections**(methodName='runTest')  
Bases: unittest.case.TestCase

TestCase class for the module *Intersections*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**  
Setup method.

**tearDown()**  
Teardown method.

**testSkeleton()**  
First test to check if the testcase is working with the testing framework.

**test\_extractFromLines()**  
Tests for method *extractFromLines*.

## pymcxray.FileFormat.Results.test\_MicroscopeParameters module

Tests for module *MicroscopeParameters*.

**class** pymcxray.FileFormat.Results.test\_MicroscopeParameters.**TestMicroscopeParameters**(methodName='runTest')  
Bases: unittest.case.TestCase

TestCase class for the module *MicroscopeParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**  
Setup method.

**tearDown()**  
Teardown method.

**testSkeleton()**  
First test to check if the testcase is working with the testing framework.

**test\_readFromLines()**  
Tests for method *readFromLines*.

pymcxray.FileFormat.Results.test\_MicroscopeParameters.**getLinesAndReference()**

## pymcxray.FileFormat.Results.test\_ModelParameters module

ModelParameters

**class** pymcxray.FileFormat.Results.test\_ModelParameters.**TestModelParameters**(methodName='runTest')  
Bases: unittest.case.TestCase

TestCase class for the module *ModelParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_readFromLines()
    Tests for method readFromLines.
```

`pymcxray.FileFormat.Results.test_ModelParameters.getLinesAndReference()`

## [pymcxray.FileFormat.Results.test\\_Phirhoz module](#)

Tests for the module *Phirhoz*.

**class** `pymcxray.FileFormat.Results.test_Phirhoz.TestPhirhoz(methodName='runTest')`  
Bases: `unittest.case.TestCase`

TestCase class for the module *Phirhoz*.

Create an instance of the class that will use the named test method when executed. Raises a `ValueError` if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_readFromLines()
    Tests for method readFromLines.
```

`pymcxray.FileFormat.Results.test_Phirhoz.getLinesAndReference(path)`

## [pymcxray.FileFormat.Results.test\\_PhirhozElement module](#)

Tests for the module *PhirhozElement*.

**class** `pymcxray.FileFormat.Results.test_PhirhozElement.TestPhirhozElement(methodName='runTest')`  
Bases: `unittest.case.TestCase`

TestCase class for the module *PhirhozElement*.

Create an instance of the class that will use the named test method when executed. Raises a `ValueError` if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_readFromLine()
    Tests for method readFromLines.
```

```
pymcxray.FileFormat.Results.test_PhirhozElement.getLineAndReference()
```

## pymcxray.FileFormat.Results.test\_PhirhozEmittedCharacteristic module

Tests for the module *PhirhozEmittedCharacteristic*.

```
class pymcxray.FileFormat.Results.test_PhirhozEmittedCharacteristic.TestPhirhozEmittedCharacteristic
    Bases: unittest.case.TestCase
```

TestCase class for the module *PhirhozEmittedCharacteristic*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.
```

```
tearDown()
    Teardown method.
```

```
testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_read()
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_PhirhozEmittedCharacteristicThinFilm module

Tests for module *PhirhozEmittedCharacteristicThinFilm*.

```
class pymcxray.FileFormat.Results.test_PhirhozEmittedCharacteristicThinFilm.TestPhirhozEmittedCharacteristicThinFilm
    Bases: unittest.case.TestCase
```

TestCase class for the module *PhirhozEmittedCharacteristicThinFilm*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.
```

```
tearDown()
    Teardown method.
```

```
testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_getIntensity()
    Tests for method getIntensity.
```

```
test_read()
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_PhirhozGenerated module

Tests for the module *PhirhozGenerated*.#

**class** pymcxray.FileFormat.Results.test\_PhirhozGenerated.**TestPhirhozGenerated**(*methodName*='runAll')  
Bases: unittest.case.TestCase

TestCase class for the module *moduleName*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_PhirhozGeneratedCharacteristic module

Tests for the module *PhirhozGeneratedCharacteristic*.

**class** pymcxray.FileFormat.Results.test\_PhirhozGeneratedCharacteristic.**TestPhirhozGeneratedCharacteristic**(*methodName*='runAll')  
Bases: unittest.case.TestCase

TestCase class for the module *PhirhozGeneratedCharacteristic*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_PhirhozGeneratedCharacteristicThinFilm module

Tests for module *PhirhozGeneratedCharacteristicThinFilm*.

**class** pymcxray.FileFormat.Results.test\_PhirhozGeneratedCharacteristicThinFilm.**TestPhirhozGeneratedCharacteristicThinFilm**(*methodName*='runAll')  
Bases: unittest.case.TestCase

TestCase class for the module *PhirhozGeneratedCharacteristicThinFilm*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_getIntensity()
    Tests for method getIntensity.

test_read()
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_PhirhozRegion module

Tests for the module *PhirhozRegion*.

```
class pymcxray.FileFormat.Results.test_PhirhozRegion.TestPhirhozRegion(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *PhirhozRegion*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_readFromLines()
    Tests for method readFromLines.
```

```
pymcxray.FileFormat.Results.test_PhirhozRegion.getLinesAndReference(path)
```

## pymcxray.FileFormat.Results.test\_RegionParameters module

Tests for the module *RegionParameters*.

```
class pymcxray.FileFormat.Results.test_RegionParameters.TestRegionParameters(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *RegionParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

## pymcxray.FileFormat.Results.test\_RegionVolume module

Tests for the module *RegionVolume*.

```
class pymcxray.FileFormat.Results.test_RegionVolume.TestRegionVolume(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *RegionVolume*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
NOtest_readFromLines()
    Tests for method readFromLines.
```

```
setUp()
    Setup method.
```

```
tearDown()
    Teardown method.
```

```
testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
pymcxray.FileFormat.Results.test_RegionVolume.getLinesAndReference()
```

## pymcxray.FileFormat.Results.test\_SimulationParameters module

Tests for the module *SimulationParameters*.

```
class pymcxray.FileFormat.Results.test_SimulationParameters.TestSimulationParameters(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *SimulationParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.
```

```
tearDown()
    Teardown method.
```

```
testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_readFromLines()
    Tests for method readFromLines.
```

```
pymcxray.FileFormat.Results.test_SimulationParameters.getLinesAndReference()
```

## pymcxray.FileFormat.Results.test\_Spectra module

Tests for the module *Spectra*.

```
class pymcxray.FileFormat.Results.test_Spectra.TestSpectra(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *Spectra*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test__extractElementHeader()
    Tests for method _extractElementHeader.

test__extractRegionHeader()
    Tests for method _extractRegionHeader.

test__read()
    Tests for method read.

test__readRegion()
    Tests for method readRegion.

test__readSpecimen()
    Tests for method readSpecimen.
```

## pymcxray.FileFormat.Results.test\_SpectraEDS module

Tests for module *SpectraEDS*.

```
class pymcxray.FileFormat.Results.test_SpectraEDS.TestSpectraEDS(methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *SpectraEDS*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test__FindTestData()
    Tests for method FindTestData.

test__extractTotalCounts()
    Tests for method _extractTotalCounts.

test__isPartialSpectraReferenceSection()
    Tests for method _isTestInputSection.

test__isRegionSpectraSection()
    Tests for method _isRegionSpectraSection.

test__isTestInputSection()
    Tests for method _isTestInputSection.
```

**test\_readPartialSpectraReferenceSection()**

Tests for method *readTestInputSection*.

**test\_readRegionSpectraSection()**

Tests for method *readRegionSpectraSection*.

**test\_readTestInputSection()**

Tests for method *readTestInputSection*.

## pymcxray.FileFormat.Results.test\_Spectrum module

Tests for the module *Spectrum*

**class pymcxray.FileFormat.Results.test\_Spectrum.TestSpectrum(methodName='runTest')**

Bases: unittest.case.TestCase

TestCase class for the module *Spectrum*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

## pymcxray.FileFormat.Results.test\_Tags module

Tests for the modules *Tags*.

**class pymcxray.FileFormat.Results.test\_Tags.TestTags(methodName='runTest')**

Bases: unittest.case.TestCase

TestCase class for the module *Tags*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**getLines()**

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_findTag()**

Tests for method *findTag*.

## pymcxray.FileFormat.Results.test\_XrayIntensities module

Tests for the module *XrayIntensities*.

**class** pymcxray.FileFormat.Results.test\_XrayIntensities.**TestXrayIntensities** (*methodName*=’runTest’)  
Bases: unittest.case.TestCase

TestCase class for the module *XrayIntensities*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_XraySimulatedSpectraRegion module

description

**class** pymcxray.FileFormat.Results.test\_XraySimulatedSpectraRegion.**TestXraySimulatedSpectraRegion** (*methodName*=’runTest’)  
Bases: unittest.case.TestCase

TestCase class for the module *moduleName*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_read()**

Tests for method *read*.

## pymcxray.FileFormat.Results.test\_XraySimulatedSpectraSpecimen module

**class** pymcxray.FileFormat.Results.test\_XraySimulatedSpectraSpecimen.**TestXraySimulatedSpectraSpecimen** (*methodName*=’runTest’)  
Bases: unittest.case.TestCase

TestCase class for the module *XraySimulatedSpectraSpecimen*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

```
tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_read()
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_XraySpectraAtomEmittedDetectedLines module

Tests for the module *XraySpectraAtomEmittedDetectedLines*.

```
class pymcxray.FileFormat.Results.test_XraySpectraAtomEmittedDetectedLines.TestXraySpectra
    Bases: unittest.case.TestCase

    TestCase class for the module XraySpectraAtomEmittedDetectedLines.

    Create an instance of the class that will use the named test method when executed. Raises a ValueError if the
    instance does not have a method with the specified name.

setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_read()
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_XraySpectraRegionEmitted module

### pymcxray.FileFormat.Results.test\_XraySpectraRegionsEmitted module

### pymcxray.FileFormat.Results.test\_XraySpectraSpecimen module

Tests for the module *XraySpectraSpecimen*.

```
class pymcxray.FileFormat.Results.test_XraySpectraSpecimen.TestXraySpectraSpecimen(methodName)
    Bases: unittest.case.TestCase

    TestCase class for the module XraySpectraSpecimen.

    Create an instance of the class that will use the named test method when executed. Raises a ValueError if the
    instance does not have a method with the specified name.

setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_read()  
    Tests for method read.
```

## pymcxray.FileFormat.Results.test\_XraySpectraSpecimenEmittedDetected module

Tests for the module *XraySpectraSpecimenEmittedDetected*.

```
class pymcxray.FileFormat.Results.test_XraySpectraSpecimenEmittedDetected.TestXraySpectraSpecimenEmittedDetected  
    Bases: unittest.case.TestCase
```

TestCase class for the module *XraySpectraSpecimenEmittedDetected*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()  
    Setup method.
```

```
tearDown()  
    Teardown method.
```

```
testSkeleton()  
    First test to check if the testcase is working with the testing framework.
```

```
test_read()  
    Tests for method read.
```

## pymcxray.FileFormat.Results.tests module

### Module contents

#### Submodules

## pymcxray.FileFormat.Element module

MCXRay element input file.

```
class pymcxray.FileFormat.Element.Element (atomicNumber=0, massFraction=1.0)  
    Bases: object
```

**atomicNumber**

**createLineOldVersion()**

**createLinesWithKey()**

**extractFromLineOldVersion(*line*)**

**extractFromLinesWithKey(*lines*)**

**massFraction**

**name**

## pymcxray.FileFormat.ExportedSpectrum module

Read and write exported spectrum from McXRay.

```
class pymcxray.FileFormat.ExportedSpectrum.ExportedSpectrum
    Bases: object

    getData()
    getSpectrumType()
    read(filepath)
```

## pymcxray.FileFormat.FileReaderWriterTools module

description

```
pymcxray.FileFormat.FileReaderWriterTools.reduceAfterDot(value)
```

## pymcxray.FileFormat.MCXRayModel module

Model type used in MCXRay.

```
class pymcxray.FileFormat.MCXRayModel.AtomCollisionModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_BROWNING = 1
    TYPE_GAUVIN = 2
    TYPE_RUTHERFORD = 0

class pymcxray.FileFormat.MCXRayModel.AtomCollisionScreeningModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_HENOC_MAURICE = 0

class pymcxray.FileFormat.MCXRayModel.AtomCrossSectionModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_BROWNING = 0
    TYPE_GAUVIN_DROUIN = 1

class pymcxray.FileFormat.MCXRayModel.AtomCrossSectionScreeningModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_HENOC_MAURICE = 0

class pymcxray.FileFormat.MCXRayModel.AtomElectronRangeModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_KANAYA_OKAYAMA = 0

class pymcxray.FileFormat.MCXRayModel.AtomEnergyLossModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_BETHE = 0

class pymcxray.FileFormat.MCXRayModel.AtomMeanIonizationPotentialModel(currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

    TYPE_JOY LUO = 0
```

```

class pymcxray.FileFormat.MCXRayModel.AtomScreeningModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPE_HENOC_MAURICE = 0

class pymcxray.FileFormat.MCXRayModel.MCXRayModel (currentModel=None)
    Bases: object

        getModel()

        setModel (modelType)
        setModelFromString (text)

class pymcxray.FileFormat.MCXRayModel.MassAbsorptionCoefficientModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPECHANTLER2005 = 3
        TYPEHEINRICH_DATA = 1
        TYPEHEINRICH_PARAMETERIZATION = 2
        TYPEHENKE = 0

class pymcxray.FileFormat.MCXRayModel.RegionEnergyLossModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPEBETHE = 1
        TYPEBETHE_JOY_LUO = 0
        TYPEBETHE_RELATIVISTIC = 2
        TYPEJOY_LUO_KGAUVIN = 3
        TYPEJOY_LUO_MONSEL = 4

class pymcxray.FileFormat.MCXRayModel.SampleEnergyLossModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPEBETHE_JOY_LUO = 0

class pymcxray.FileFormat.MCXRayModel.SpectrumInterpolationModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPECOPY = 0
        TYPELINEAR = 1
        TYPELINEAR_DOUBLE = 2
        TYPESPLINE = 3
        TYPESPLINE_BATCH = 4
        TYPESPLINE_POINT = 5

class pymcxray.FileFormat.MCXRayModel.XRayCSBremsstrahlungModel (currentModel=None)
    Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

        TYPEBETHE_HEITLER = 0
        TYPERDING = 2
        TYPEGAUVIN = 3
        TYPEKIRKPATRICK_WIEDMAN = 1

```

```
class pymcxray.FileFormat.MCXRayModel.XRayCSCharacteristicModel (currentModel=None)
Bases: pymcxray.FileFormat.MCXRayModel.MCXRayModel

TYPE_BOTE2009 = 1
TYPE_CASTANI1982 = 0
```

## pymcxray.FileFormat.MicroscopeParameters module

MCXRay microscope parameters input file.

```
class pymcxray.FileFormat.MicroscopeParameters.MicroscopeParameters
Bases: object

beamCurrent_A
beamDiameter_A
beamEnergy_keV
beamPositionX_A
beamPositionY_A
beamStandardDeviation_A
beamTilt_deg
defaultValues()
detectorAzimuthalAngle_deg
detectorBFHigh_rad
detectorBFLow_rad
detectorChannelWidth_eV
detectorCrystalAtomSymbol
detectorCrystalDistance_cm
detectorCrystalRadius_cm
detectorCrystalThickness_cm
detectorDFHigh_rad
detectorDFLow_rad
detectorDeadLayer_A
detectorDiffusionLength_A
detectorHAADFHigh_rad
detectorHAADFLow_rad
detectorNoise_eV
detectorPitch_deg
detectorSurfaceQuality
detectorTOA_deg
read(filepath)
```

```
time_s
version
write (filepath)
```

## pymcxray.FileFormat.Models module

MCXRay models file.

```
class pymcxray.FileFormat.Models.Models
    Bases: object

    getModelList ()
    modelAtomCollision
    modelAtomCrossSection
    modelAtomMac
    modelSampleEnergyLoss
    modelXrayBremsstrahlung
    modelXrayCharacteristic
    read (filepath)
    version
    write (filepath)
```

## pymcxray.FileFormat.Region module

MCXRay region input file.

```
class pymcxray.FileFormat.Region.Region
    Bases: object

    clear ()
    createLinesWithVersion ()
    createLinesWithoutVersion ()
    extractFromLinesWithVersion (lines)
    extractFromLinesWithoutVersion (lines)
```

## pymcxray.FileFormat.RegionDimensions module

MCXRay region dimensions input file.

```
class pymcxray.FileFormat.RegionDimensions.RegionDimensions (parameters=None)
    Bases: object

    createLineOldVersion ()
    createLineWithKey ()
    extractFromLineOldVersion (line)
```

```
extractFromLinesWithKey (line)
getLineFormat ()

class pymcxray.FileFormat.RegionDimensions.RegionDimensionsBox (parameters=None)
    Bases: pymcxray.FileFormat.RegionDimensions.RegionDimensions

    getLineFormat ()

    maximumX
    maximumY
    maximumZ
    minimumX
    minimumY
    minimumZ

class pymcxray.FileFormat.RegionDimensions.RegionDimensionsCylinder (parameters=None)
    Bases: pymcxray.FileFormat.RegionDimensions.RegionDimensionsSphere

    directionX
    directionY
    directionZ
    getLineFormat ()
    length

class pymcxray.FileFormat.RegionDimensions.RegionDimensionsSphere (parameters=None)
    Bases: pymcxray.FileFormat.RegionDimensions.RegionDimensions

    getLineFormat ()

    positionX
    positionY
    positionZ
    radius

pymcxray.FileFormat.RegionDimensions.createRegionDimensions (regionType)
```

## pymcxray.FileFormat.RegionType module

MCXRay region type input file.

## pymcxray.FileFormat.ResultsParameters module

MCXRay ResultsParameters input file.

```
class pymcxray.FileFormat.ResultsParameters.ResultsParameters
    Bases: object

    defaultValues ()
    isComputeXrayBremsstrahlung
    isComputeXrayCharacteristic
```

```
isComputeXrayPhirhoz
isComputeXraySimulatedSpectrum
read(filepath)
version
write(filepath)
```

## pymcxray.FileFormat.SimulationInputs module

MCXRay simulation inputs file.

```
class pymcxray.FileFormat.SimulationInputs.SimulationInputs
Bases: object
getExtension(key)
mapFilename
microscopeFilename
modelFilename
read(filepath)
resultParametersFilename
simulationParametersFilename
snrFilename
specimenFilename
title
version
write(filepath)
```

## pymcxray.FileFormat.SimulationParameters module

MCXRay simulation parameters input file.

```
class pymcxray.FileFormat.SimulationParameters.SimulationParameters
Bases: object
baseFilename
defaultValues()
elasticCrossSectionScalingFactor
energyChannelWidth_eV
energyLossScalingFactor
numberChannels
numberElectrons
numberFilmsX
numberFilmsY
```

```
numberFilmsZ
numberPhotons
numberWindows
read(filepath)
spectrumInterpolationModel
version
voxelSimplification
write(filepath)
```

## pymcxray.FileFormat.SnrParameters module

MCXRay Snr input file.

```
class pymcxray.FileFormat.SnrParameters.SnrParameters
Bases: object
backgroundEnergyWindowSize
defaultValues()
energyEnd_keV
energyStart_keV
numberEnergySteps
read(filepath)
snrType
spectrumEnergyWindowSize
write(filepath)
```

## pymcxray.FileFormat.Specimen module

MCXRay specimen input file.

```
class pymcxray.FileFormat.Specimen.Specimen
Bases: object
clear()
name
numberRegions
read(filepath)
regions
version
write(filepath)
```

## pymcxray.FileFormat.Version module

MCXray version information.

```
class pymcxray.FileFormat.Version.Version(major, minor, revision)
    Bases: object

    fromString(versionString)

    key = 'Version'

    major

    minor

    readFromFile(filepath)

    revision

    toString()

    writeLine(outputFile)
```

## pymcxray.FileFormat.testUtilities module

Utilities used for testing this package.

```
pymcxray.FileFormat.testUtilities.createTempDataPath(path)
pymcxray.FileFormat.testUtilities.getSimulationTitles()
pymcxray.FileFormat.testUtilities.removeTempDataPath(path)
```

## pymcxray.FileFormat.test\_Element module

Tests for module *Element*.

```
class pymcxray.FileFormat.test_Element.TestElement(methodName='runTest')
    Bases: unittest.case.TestCase
```

TestCase class for the module *Element*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Setup method.
```

```
tearDown()
    Teardown method.
```

```
testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

```
test_createLineOldVersion()
    Tests for method createLineOldVersion.
```

```
test_createLineWithKey()
    Tests for method createLineOldVersion.
```

```
test_extractFromLineOldVersion()
    Tests for method extractFromLineOldVersion.
```

```
test_extractFromLineWithKey()  
    Tests for method extractFromLinesWithKey.
```

## pymcxray.FileFormat.test\_ExportedSpectrum module

Tests for module *ExportedSpectrum*.

```
class pymcxray.FileFormat.test_ExportedSpectrum.TestExportedSpectrum(methodName='runTest')  
    Bases: unittest.case.TestCase
```

TestCase class for the module *moduleName*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()  
    Setup method.
```

```
tearDown()  
    Teardown method.
```

```
testSkeleton()  
    First test to check if the testcase is working with the testing framework.
```

```
test_read()  
    Tests for method read.
```

## pymcxray.FileFormat.test\_FileReaderWriterTools module

Tests for the module *FileReaderWriterTools*.

```
class pymcxray.FileFormat.test_FileReaderWriterTools.TestFileReaderWriterTools(methodName='runTest')  
    Bases: unittest.case.TestCase
```

TestCase class for the module *FileReaderWriterTools*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()  
    Setup method.
```

```
tearDown()  
    Teardown method.
```

```
testSkeleton()  
    First test to check if the testcase is working with the testing framework.
```

```
test_reduceAfterDot()  
    Tests for method reduceAfterDot.
```

## pymcxray.FileFormat.test\_MCXRayModel module

Tests for module *MCXRayModel*.

```
class pymcxray.FileFormat.test_MCXRayModel.TestMCXRayModel(methodName='runTest')  
    Bases: unittest.case.TestCase
```

TestCase class for the module *MCXRayModel*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_MCXRayModel()**

Tests for method *MCXRayModel*.

**pymcxray.FileFormat.test\_MicroscopeParameters module****pymcxray.FileFormat.test\_Models module****pymcxray.FileFormat.test\_Region module**

Tests for module *Region*.

**class pymcxray.FileFormat.test\_Region.TestRegion(methodName='runTest')**

Bases: unittest.case.TestCase

TestCase class for the module *moduleName*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**getTestRegionLinesWithVersion(title, userMassDensity=True)****getTestRegionLinesWithoutVersion(title, userMassDensity=True)****setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_createLinesWithVersion()**

Tests for method *createLinesWithVersion*.

**test\_createLinesWithoutVersion()**

Tests for method *createLinesWithoutVersion*.

**test\_extractFromLinesWithVersion()**

Tests for method *read*.

**test\_extractFromLinesWithoutVersion()**

Tests for method *read*.

**pymcxray.FileFormat.test\_RegionDimensions module**

Tests for module *RegionDimensions*.

```
class pymcxray.FileFormat.test_RegionDimensions.TestRegionDimensions (methodName='runTest')
Bases: unittest.case.TestCase

TestCase class for the module RegionDimensions.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the
instance does not have a method with the specified name.

setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.

test_RegionDimensionsBox_createLineOldVersion()
    Tests for class RegionDimensionsBox.

test_RegionDimensionsBox_createLineWithKey()
    Tests for class RegionDimensionsBox.

test_RegionDimensionsBox_extractFromLineOldVersion()
    Tests for class RegionDimensionsBox.

test_RegionDimensionsBox_extractFromLinesWithKey()
    Tests for class RegionDimensionsBox.

test_RegionDimensionsCylinder_createLineOldVersion()
    Tests for class RegionDimensionsCylinder.

test_RegionDimensionsCylinder_createLineWithKey()
    Tests for class RegionDimensionsCylinder.

test_RegionDimensionsCylinder_extractFromLineOldVersion()
    Tests for class RegionDimensionsCylinder.

test_RegionDimensionsCylinder_extractFromLinesWithKey()
    Tests for class RegionDimensionsCylinder.

test_RegionDimensionsSphere_createLineOldVersion()
    Tests for class RegionDimensionsSphere.

test_RegionDimensionsSphere_createLineWithKey()
    Tests for class RegionDimensionsSphere.

test_RegionDimensionsSphere_extractFromLineOldVersion()
    Tests for class RegionDimensionsSphere.

test_RegionDimensionsSphere_extractFromLinesWithKey()
    Tests for class RegionDimensionsSphere.

test_createRegionDimensions()
    Tests for method createRegionDimensions.
```

## pymcxray.FileFormat.test\_RegionType module

Test for module *RegionType*.

```
class pymcxray.FileFormat.test_RegionType.TestRegionType (methodName='runTest')
Bases: unittest.case.TestCase
```

TestCase class for the module *RegionType*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_Constants()**

Tests for method *Constants*.

**pymcxray.FileFormat.test\_ResultsParameters module****pymcxray.FileFormat.test\_SimulationInputs module****pymcxray.FileFormat.test\_SimulationParameters module****pymcxray.FileFormat.test\_SnrParameters module**

Tests for the module *SnrParameters*.

**class** pymcxray.FileFormat.test\_SnrParameters.**TestSnrParameters**(*methodName='runTest'*)  
Bases: unittest.case.TestCase

TestCase class for the module *SnrParameters*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**getSnrParametersReference(*title*)****setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_\_createKeys()**

Tests for method *\_createKeys*.

**test\_read()**

Tests for method *read*.

**test\_write()**

Tests for method *write*.

**pymcxray.FileFormat.test\_Specimen module**

## pymcxray.FileFormat.test\_Version module

Tests for module *Version*.

**class** pymcxray.FileFormat.test\_Version.**TestVersion**(*methodName*='runTest')

Bases: unittest.case.TestCase

TestCase class for the module *Version*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

**test\_VersionConstants()**

Tests for method *VersionConstants*.

**test\_comparison()**

Test comparison operation on Version class.

**test\_fromString()**

Tests for method *fromString*.

**test\_readFromFile()**

Tests for method *readFromFile*.

**test\_readFromFile\_BadFile()**

Tests for method *readFromFile*.

**test\_toString()**

Tests for method *toString*.

**test\_writeLine()**

Tests for method *writeLine*.

## pymcxray.FileFormat.tests module

### Module contents

#### pymcxray.serialization package

##### Submodules

#### pymcxray.serialization.SerializationH5py module

**class** pymcxray.serialization.SerializationH5py.**SerializationNumpy**(*filename*=None,

*ver-*  
*bose*=True)

Bases: pymcxray.serialization.\_Serialization.\_Serialization

**load()**

```
    save (serializedData)
```

## pymcxray.serialization.SerializationNumpy module

```
class pymcxray.serialization.SerializationNumpy.SerializationNumpy (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization._Serialization._Serialization  
  
    load()  
  
    save (serializedData)  
  
class pymcxray.serialization.SerializationNumpy.SerializationNumpyNPY (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization.SerializationNumpy.SerializationNumpy  
  
    load()  
  
    save (data)  
  
class pymcxray.serialization.SerializationNumpy.SerializationNumpyNPZ (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization.SerializationNumpy.SerializationNumpy  
  
    load()  
  
    save (data)  
  
class pymcxray.serialization.SerializationNumpy.SerializationNumpyTxt (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization.SerializationNumpy.SerializationNumpy  
  
    load()  
  
    save (data)  
  
class pymcxray.serialization.SerializationNumpy.SerializationNumpyTxtGz (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization.SerializationNumpy.SerializationNumpyTxt
```

## pymcxray.serialization.SerializationPickle module

```
class pymcxray.serialization.SerializationPickle.SerializationPickle (filename=None,  
                                         ver-  
                                         bosse=True)  
Bases: pymcxray.serialization._Serialization._Serialization  
  
    KEY_FILE_VERSION = 'fileVersion'  
    KEY_SERIALIZED_DATA = 'serializedData'  
  
    load()  
  
    save (serializedData)
```

## pymcxray.serialization.test\_Serialization module

### pymcxray.serialization.test\_SerializationH5py module

```
class pymcxray.serialization.test_SerializationH5py.TestSerializationH5py (methodName='runTest')  
Bases: unittest.case.TestCase
```

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Hook method for setting up the test fixture before exercising it.

**tearDown()**

Hook method for deconstructing the test fixture after testing it.

**testSkeleton()**

### pymcxray.serialization.test\_SerializationNumpy module

```
class pymcxray.serialization.test_SerializationNumpy.TestSerializationNumpy (methodName='runTest')  
Bases: unittest.case.TestCase
```

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Hook method for setting up the test fixture before exercising it.

**tearDown()**

Hook method for deconstructing the test fixture after testing it.

**testSkeleton()**

**test\_loadSaveSerializationNumpy()**

**test\_loadSaveSerializationNumpyNPY()**

**test\_loadSaveSerializationNumpyNPZ()**

**test\_loadSaveSerializationNumpyTxt()**

**test\_loadSaveSerializationNumpyTxtGz()**

### pymcxray.serialization.test\_SerializationPickle module

```
class pymcxray.serialization.test_SerializationPickle.TestSerialization (methodName='runTest')  
Bases: unittest.case.TestCase
```

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Hook method for setting up the test fixture before exercising it.

**tearDown()**

Hook method for deconstructing the test fixture after testing it.

**testSkeleton()**

```
test_loadSave()
test_version()
```

## pymcxray.serialization.tests module

Regression testing for the project.

### Module contents

#### pymcxray.tests package

##### Submodules

#### pymcxray.tests.test\_pymcxray module

##### test\_pymcxray

Tests for *pymcxray* module.

```
class pymcxray.tests.test_pymcxray.TestPymcxray(methodName='runTest')
Bases: unittest.case.TestCase
```

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()
    Hook method for setting up the test fixture before exercising it.

tearDown()
    Hook method for deconstructing the test fixture after testing it.

test_000_something()
```

### Module contents

## 8.1.2 Submodules

### 8.1.3 pymcxray.AnalyzeNumberBackgroundWindows module

Analyze the number of background windows on the x-ray spectrum.

```
class pymcxray.AnalyzeNumberBackgroundWindows.AnalyzeNumberBackgroundWindows
Bases: object

plotData()
plotDifference()
readData()

pymcxray.AnalyzeNumberBackgroundWindows.run()
```

## 8.1.4 pymcxray.AtomData module

MCXRay atom data.

```
pymcxray.AtomData.getAtomSymbol(atomic_number)
pymcxray.AtomData.getAtomicNumber(symbol)
pymcxray.AtomData.getIonizationEnergy_keV(shell, element)
pymcxray.AtomData.getMassDensity_g_cm3(symbol)
pymcxray.AtomData.getShellList()
pymcxray.AtomData.getXRayEnergy_keV(line, element)
pymcxray.AtomData.get_atomic_weight_g_mol(atomic_number)
pymcxray.AtomData.run()
```

## 8.1.5 pymcxray.BatchFile module

MCXRay batch file creator.

```
class pymcxray.BatchFile.BatchFile(name, numberFiles=1)
Bases: object
    addSimulationName(simulationFilename)
    removePreviousFiles(path)
    write(path)
```

## 8.1.6 pymcxray.BatchFileConsole module

MCXRay console batch file creator.

```
class pymcxray.BatchFileConsole.BatchFileConsole(name,      programName,      number-
Files=1)
Bases: object
```

The batch file is responsible to create the simulation structure with a copy of mcxray program.

One important parameter to set is the *numberFiles*, this is the number of batch files generated and that can be run in parallel. For maximum efficiency it should be set as the number of logical processors minus 1 or 2. For example, on a computer with 12 logical processors, the *numberFiles* should be set at 10.

### Parameters

- **name** (*str*) – Basename used for the batch files
- **programName** (*str*) – Name of the executable to add in the batch file
- **numberFiles** (*int*) – Number of batch files to generate and possibly to run in parallel

**addSimulationName** (*simulationFilename*)

Add a simulation in the simulation list.

**Parameters** **simulationFilename** (*str*) – File path of the simulation added

**write** (*path*)

Write the batch files for all simulations in the simulation list.

**Parameters** **path** (*str*) – Path where the batch files are written.

## 8.1.7 pymcxray.ComparisonModels module

Comparison of the models used by MCXray.

```
class pymcxray.ComparisonModels.ComparisonModels(dataPath)
Bases: object

graphicsEnergyLoss()
graphicsIonizationCrossSection()
graphicsXrayCrossSectionBremstrahlung()
graphicsXrayMassAbsorptionCoefficient()

pymcxray.ComparisonModels.runVersion1_2_3()
pymcxray.ComparisonModels.runVersion1_4_0()
pymcxray.ComparisonModels.runVersion1_4_1()
```

## 8.1.8 pymcxray.DebugSimulatedSpectrum module

Debug the simulated spectrum implementation in mcxray.

```
class pymcxray.DebugSimulatedSpectrum.DebugSimulatedSpectrum
Bases: object

runRegion()
runSpecimen()

pymcxray.DebugSimulatedSpectrum.run()
```

## 8.1.9 pymcxray.ElementProperties module

```
pymcxray.ElementProperties.computeAtomicDensity_atom_cm3(massDensity_g_cm3,
atomicMass_g_mol)
```

Compute the atomic density.

$$n_i = \frac{N_A \rho_i}{A_i}$$

where

- $n_i$  is the atomic density in atoms/ $cm^3$
- $N_A$  is the Avogadro number in atoms/ $mole$
- $\rho_i$  is the mass density in  $g/cm^3$
- $A_i$  is the atomic mass in  $g/mole$

### Parameters

- **massDensity\_g\_cm3** (*float*) –
- **atomicMass\_g\_mol** (*float*) –

For element H to Lr (1-103). From: CASINO source code. DOS version.

**Todo:** Add units.

For element H to Sg (1-106).

Unit g/mole.

From: Tableau periodique des elements, Sargent-Welch scientifique Canada Limitee.

For element H to Lr (1–103). From: CASINO source code, DOS version.

**Todo:** Add units.

pymcxray.ElementProperties.g\_massDensity\_g\_cm3 = [0.0899, 0.1787, 0.53, 1.85, 2.34, 2.62, ...  
Mass density of element in atomic number order.

For element H to Cm (1-96).

In  $q/cm^3$ .

From: Tableau periodique des elements, Sargent-Welch scientifique Canada Limitee.

**Note:** Element Z = 85 and 87 set to 1 for the calculation.

```
pymcxray.ElementProperties.g_plasmonEnergy = [15.0, 15.0, 7.1, 18.7, 22.7, 15.0, 15.0, 15.0, 15.0]
```

Plasmon energy of element in atomic number order.

For element H to Lr (1-103). From: CASINO source code, DOS version.

**Todo:** Add units.

```
pymcxray.ElementProperties.getAtomicMass_g_mol(atomicNumber)
```

```
pymcxray.ElementProperties.getAtomicNumber(atomicNumber=None, name=None, symbol=None)
```

```
pymcxray.ElementProperties.getAtomicNumberByName(name)
```

```
pymcxray.ElementProperties.getAtomicNumberBySymbol(symbol)
```

```
pymcxray.ElementProperties.getFermiEnergy_eV(atomicNumber)
```

```
pymcxray.ElementProperties.getKFermi_eV(atomicNumber)
```

`pymcxray.ElementProperties.getKRatioCorrection(atomicNumber)`

Get the constant k ratio correction needed by the mean ionization potential from the atomic number.

Parameters **atomic\_number** (*int*) – Atomic number

---

```
pymcxray.ElementProperties.getKRatioCorrectionMonsel(atomicNumber, workFunction_keV)
    /// K value as defined by Monsel. /// Used in DE/DS calculation. Casino uses K Gauvin, but for low energy, ///
    JR Lowney says that this one is more appropriate (and by experience, // it is effectively better for the secondary
    yield). /// <p> NOTE : Depends on J (ionisation potential). So it must already be calculated before. /// @param
    element Element for whom we want to calculate the K value. /// @return The K value of the element passed in
    argument
```

`pymcxray.ElementProperties.getMassDensity_g_cm3(atomicNumber)`

`pymcxray.ElementProperties.getMeanIonizationEnergy_eV(atomic_number)`

Get the mean ionization potential from the atomic number.

In eV.

**Parameters** `atomic_number` (`int`) – Atomic number

`pymcxray.ElementProperties.getName(atomicNumber)`

`pymcxray.ElementProperties.getPlasmonEnergy_eV(atomicNumber)`

`pymcxray.ElementProperties.getSymbol(atomicNumber)`

`pymcxray.ElementProperties.run()`

`pymcxray.ElementProperties.runAtomicNumberSymbol()`

## 8.1.10 pymcxray.Simulation module

MCXRay simulation parameters.

`class pymcxray.Simulation.Layer(elements, thickness_nm, mass_density_g_cm3=None)`  
Bases: object

`class pymcxray.Simulation.Simulation(overwrite=True)`  
Bases: object

`basename`

`beamDiameter_nm`

`beamPosition_nm`

`beamTilt_deg`

`createSimulationFiles(path, simulationPath, hdf5_group)`

`current_A`

`detectorAzimuthalAngle_deg`

`detectorChannelWidth_eV`

`detectorCrystalDistance_cm`

`detectorCrystalRadius_cm`

`detectorCrystalThickness_cm`

`detectorNoise_eV`

`elasticCrossSectionScalingFactor`

`energyLossScalingFactor`

`energy_keV`

```
filename
generateBaseFilename()
getFilenameSuffixes()
getParameters()
getProgramVersionFilepath(simulationPath)
isDone(simulationPath, hdf5_group=None)
modelXrayBremsstrahlung
name
numberContinuumWindows
numberElectrons
numberEnergyWindows
numberLayersX
numberLayersY
numberLayersZ
numberPhotons
removeInputsFiles()
resultsBasename
setParameters(parameters)
solidAngle_sr
spectrumInterpolationModel
takeOffAngle_deg
time_s

pymcxray.Simulation.computeWeightFraction(atomicNumberRef, atomicWeights)
pymcxray.Simulation.createAlloyBoxInSubstrate(elementsParticle, elementsSubstrate, box-
Parameters_nm)
pymcxray.Simulation.createAlloyBoxInThinFilm(elementsParticle, atomicNumberSubstrate,
boxParameters_nm, filmThickness_nm)
pymcxray.Simulation.createAlloyBoxInVaccuum(elements, boxParameters_nm)
pymcxray.Simulation.createAlloyBulkSample(elements, sampleName=None)
pymcxray.Simulation.createAlloyFilmOverSubstrate(film_elements, substrate_elements,
film_thickness_nm=10.0,
film_mass_density_g_cm3=None,
substrate_mass_density_g_cm3=None)
pymcxray.Simulation.createAlloyMultiVerticalLayer(elementsLayers, layerWidths_nm)
pymcxray.Simulation.createAlloyParticleInSubstrate(elementsParticle, atomicNum-
berSubstrate, particleRadius_nm,
particlePositionZ_nm=None)
```

```
pymcxray.Simulation.createAlloyParticleInThinFilm(elementsParticle, atomicNumberSubstrate, particleRadius_nm, filmThickness_nm, particlePositionZ_nm=None)
pymcxray.Simulation.createAlloyThinFilm(elements, filmThickness_nm)
pymcxray.Simulation.createAlloyThinFilm2(elements, filmThickness_nm)
pymcxray.Simulation.createBoxFeatureInSubstrate(feature_elements, substrate_elements, depth_nm, width_nm)
pymcxray.Simulation.createFilmInSubstrate(atomicNumberFilm, atomicNumberSubstrate, filmThickness_nm, filmTopPositionZ_nm)
pymcxray.Simulation.createFilmOverSubstrate(atomicNumberFilm, atomicNumberSubstrate, filmThickness_nm=10.0)
pymcxray.Simulation.createParticleInSubstrate(atomicNumberParticle, atomicNumberSubstrate, particleRadius_nm, particlePositionZ_nm=None)
pymcxray.Simulation.createParticleOnFilm(atomicNumberParticle, atomicNumberSubstrate, particleDiameter_nm, filmThickness_nm)
pymcxray.Simulation.createParticleOnSubstrate(atomicNumberParticle, atomicNumberSubstrate, particleDiameter_nm)
pymcxray.Simulation.createPhirhozSpecimens(atomicNumberTracer, atomicNumberMatrix, tracerThickness_nm, maximumThickness_nm)
pymcxray.Simulation.createPureBulkSample(atomic_number)
pymcxray.Simulation.create_cnt_sample(body_elements, cnt_length_nm=1000.0, cnt_outside_diameter_nm=100.0, cnt_inside_diameter_nm=50.0, particle_diameter_nm=5.0)
pymcxray.Simulation.create_multi_horizontal_layer(substrate_elements, layers, substrate_mass_density_g_cm3=None)
```

Create a horizontal multi layer sample.

The substrate is the first region created. The other region are each of the element in the *layers*.

#### Parameters

- **substrate\_elements** – list of atomic number and weight fraction pair for the composition of the substrate
- **layers** –
- **substrate\_mass\_density\_g\_cm3** –

#### Returns

```
pymcxray.Simulation.create_weight_fractions(weight_fraction_step, number_elements)
pymcxray.Simulation.create_weight_fractions_trace(weight_fraction_step, number_elements, min_weight_fraction, max_weight_fraction)
```

### 8.1.11 pymcxray.SimulationsParameters module

Simulations parameters

```
class pymcxray.SimulationsParameters.SimulationsParameters
Bases: dict

    addCompute (parameterKey)
    addFixed (parameterKey, value)
    addVaried (parameterKey, values)
    computeNumberXrays (experiment)
    computeNumberXraysFilepath
    fixedParameters
    getAllSimulationParameters ()
    getVariedParameterLabels ()
    variedParameters

class pymcxray.SimulationsParameters.SimulationsParametersFixed
Bases: object

    addExperiment (experiment)
    getAllSimulationParameters ()
```

## 8.1.12 pymcxray.Testings module

## 8.1.13 pymcxray.mcxray module

Base module to create and analuyze MCXRay simulations.

## 8.1.14 pymcxray.multipleloop module

This module provides a tool for handling computer experiments with of a set of input parameters, where each input parameter is varied in a prescribed fashion.

In short, the parameters are held in a dictionary where the keys are the names of the parameters and the values are the numerical, string or other values of the parameters. The value can take on multiple values: e.g., an integer parameter ‘a’ can have values -1, 1 and 10. Similarly, a string parameter ‘method’ can have values ‘Newton’ and ‘Bisection’. The module will generate all combination of all parameters and values, which in the mentioned example will be (-1, ‘Newton’), (1, ‘Newton’), (10, ‘Newton’), (-1, ‘Bisection’), (1, ‘Bisection’), and (10, ‘Bisection’). Particular combination of values can easily be removed.

The usage and implementation of the module are documented in the book “Python Scripting for Computational Science” (H. P. Langtangen, Springer, 2009), Chapter 12.1.

```
pymcxray.multipleloop.combine (prm_values)
```

Compute the combination of all parameter values in the prm\_values (nested) list. Main function in this module.

```
param prm_values: nested list (parameter_name, list_of_parameter_values) or dictionary
prm_values[parameter_name] = list_of_parameter_values. return: (all, names, varied)
where
```

- all contains all combinations (experiments) all[i] is the list of individual parameter values in experiment no i
- names contains a list of all parameter names

- varied holds a list of parameter names that are varied (i.e. where there is more than one value of the parameter, the rest of the parameters have fixed values)

Code example:

```
>>> dx = array([1.0/2**k for k in range(2,5)])
>>> dt = 3*dx; dt = dt[:-1]
>>> p = {'dx': dx, 'dt': dt}
>>> p
{'dt': [ 0.75 , 0.375,], 'dx': [ 0.25 , 0.125 , 0.0625,]}
>>> all, names, varied = combine(p)
>>> all
[[0.75, 0.25], [0.375, 0.25], [0.75, 0.125], [0.375, 0.125],
 [0.75, 0.0625], [0.375, 0.0625]]
```

## 8.1.15 pymcxray.pymcxray module

### 8.1.16 pymcxray.test\_AtomData module

Tests for the module *AtomData*.

**class** pymcxray.test\_AtomData.**TestAtomData** (*methodName='runTest'*)  
Bases: unittest.case.TestCase

TestCase class for the module *AtomData*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

**testConstants()**

First test to check if the testcase is working with the testing framework.

**testSkeleton()**

First test to check if the testcase is working with the testing framework.

## 8.1.17 pymcxray.test\_BatchFileConsole module

Tests for the module *BatchFileConsole*.

**class** pymcxray.test\_BatchFileConsole.**TestBatchFileConsole** (*methodName='runTest'*)  
Bases: unittest.case.TestCase

TestCase class for the module *BatchFileConsole*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

**setUp()**

Setup method.

**tearDown()**

Teardown method.

```
testSkeleton()  
First test to check if the testcase is working with the testing framework.
```

## 8.1.18 pymcxray.test\_ComparisonModels module

Tests for the module *ComparisonModels*.

```
class pymcxray.test_ComparisonModels.TestComparisonModels(methodName='runTest')  
Bases: unittest.case.TestCase
```

TestCase class for the module *ComparisonModels*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()  
Setup method.
```

```
tearDown()  
Teardown method.
```

```
testSkeleton()  
First test to check if the testcase is working with the testing framework.
```

## 8.1.19 pymcxray.test\_Simulation module

Tests for the module *Simulation*.

```
class pymcxray.test_Simulation.TestSimulation(methodName='runTest')  
Bases: unittest.case.TestCase
```

TestCase class for the module *Simulation*.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the instance does not have a method with the specified name.

```
setUp()  
Setup method.
```

```
tearDown()  
Teardown method.
```

```
testSkeleton()  
First test to check if the testcase is working with the testing framework.
```

```
test_create_multi_horizontal_layer()  
Test the create_multi_horizontal_layer method.
```

```
test_create_weight_fractions()
```

```
test_create_weight_fractions_trace()
```

```
test_create_weight_fractions_trace_speed_test_three_elements()
```

```
test_create_weight_fractions_trace_speed_test_two_elements()
```

## 8.1.20 pymcxray.test\_SimulationsParameters module

Tests for the module *SimulationsParameters*.

```
class pymcxray.test_SimulationsParameters.TestSimulationsParameters (methodName='runTest')
Bases: unittest.case.TestCase

TestCase class for the module SimulationsParameters.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the
instance does not have a method with the specified name.

setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

## 8.1.21 pymcxray.test\_mcxray module

Tests for the module *mcxray*.

```
class pymcxray.test_mcxray.Testmcxray (methodName='runTest')
Bases: unittest.case.TestCase

TestCase class for the module mcxray.

Create an instance of the class that will use the named test method when executed. Raises a ValueError if the
instance does not have a method with the specified name.

setUp()
    Setup method.

tearDown()
    Teardown method.

testSkeleton()
    First test to check if the testcase is working with the testing framework.
```

## 8.1.22 pymcxray.tests module

### 8.1.23 Module contents

`pymcxray.create_path(path)`

Create a path from the input string if does not exists.

Does not try to distinct between file and directory in the input string. `path = "dir1/filename.ext" => "dir1/filename.ext"` where the new directory “filename.ext” is created.

**Parameters** `path (str)` – The path input string.

**Returns** The path with the path separator at the end

**Return type** str

```
pymcxray.find_all_files(root, patterns='*', ignore_path_patterns='', ignore_name_patterns='',
single_level=False, yield_folders=False)
```

Find all files in a root folder.

From Python Cookbook section 2.16 pages 88–90

**Parameters**

- **root** –
- **patterns** –
- **ignore\_path\_patterns** –
- **ignore\_name\_patterns** –
- **single\_level** –
- **yield\_folders** –

**Returns**

`pymcxray.get_current_module_path(module_path, relative_path=’’)`

Extract the current module path and combine it with the relative path and return it.

**Parameters**

- **module\_path** (*str*) – Pass the `__FILE__` python keyword for this parameter
- **relative\_path** (*str*) – The relative path to combine with the module path

**Returns** The path obtained when combine the module path and relative path

**Return type** str

`pymcxray.get_mcxray_archive_name(configuration_file_path, default=None)`

Read the MCXRay archive name in the configuration file. This option allows to choose which version of MCXRay to use for the simulations.

The configuration file need to have this entry in the section [Paths]:

```
[Paths]
mcxrayArchiveName=2016-04-11_11h41m28s_MCXRay_v1.6.6.zip
```

**Parameters**

- **configuration\_file\_path** (*str*) – The file path of the configuration file
- **default** (*str*) – Default value to use if the entry is not found

**Returns** The MCXRay archive name

**Return type** str

`pymcxray.get_mcxray_archive_path(configuration_file_path, relative_path=’’)`

Read the MCXRay archive path in the configuration file.

The configuration file need to have this entry in the section [Paths]:

```
[Paths]
mcxrayArchivePath=D:\Dropbox\hdemers\professional\softwareRelease\mcxray
```

**Parameters**

- **configuration\_file\_path** (*str*) – The file path of the configuration file
- **relative\_path** (*str*) – Relative path to add to the path read in the configuration file

**Returns** Path where the mcxray archive can be found.

**Return type** str

`pymcxray.get_mcxtay_program_name(configuration_file_path, default=None)`

Read the MCXRay program name in the configuration file.

This option specify which executable to use in the script. The *console\_mcxtay\_x64.exe* should be OK for most situation. If you have a 32-bit system, you have to use *console\_mcxtay.exe* (32-bit version).

The configuration file need to have this entry in the section [Paths]:

```
[Paths]
mcxtayProgramName=console_mcxtay_x64.exe
```

#### Parameters

- **configuration\_file\_path** (*str*) – The fule path of the configuration file
- **default** (*str*) – Default value to use if the entry is not found

**Returns** The MCXRay program name

**Return type** str

`pymcxray.get_mcxtay_program_path(configuration_file_path, relative_path="")`

Read the MCXRay program path in the configuration file.

The configuration file need to have this entry in the section [Paths]:

```
[Paths]
mcxtayProgramPath=C:\hdemers\codings\devcasino
```

#### Parameters

- **configuration\_file\_path** (*str*) – The file path of the configuration file
- **relative\_path** (*str*) – Relative path to add to the path read in the configuration file

**Returns** Path where the mcxtay program can be found.

**Return type** str

Deprecated since version 0.1.

#### See also:

function `pymcxray.get_mcxtay_archive_path()`

`pymcxray.get_results_mcgill_path(configuration_file_path, relative_path="")`

Read the results path for McGill in the configuration file. The results path read in the configuration file is combine with the *relative\_path* and return.

The configuration file need to have this entry in the section [Paths]:

```
[Paths]
resultsMcGillPath=D:\Dropbox\hdemers\professional\results\simulations
```

#### Parameters

- **configuration\_file\_path** (*str*) – The file path of the configuration file
- **relative\_path** (*str*) – Relative path to add to the path read in the configuration file

**Returns** Path where the simulation input and results will be written

**Return type** str

```
pymcxray.read_value_from_configuration_file(configuration_file, section_name, key_name,  
                                              default=None)
```

Read a value from an entry in a section from a configuration file.

**Parameters**

- **configuration\_file** (*str*) – The file path of the configuration file
- **section\_name** (*str*) – Name of the section
- **key\_name** (*str*) – Name of the entry to read
- **default** – Default value of the entry if not found

**Returns** The value read or default value

**Return type** str

# CHAPTER 9

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