
ElikoPy

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ElikoPy is Python library aiming at easing the processing of diffusion imaging for microstructural analysis. This Python library is based on

- DIPY, a python library for the analysis of MR diffusion imaging.
- Microstructure fingerprinting, a python library doing estimation of white matter microstructural properties from a dictionary of Monte Carlo diffusion MRI fingerprints.
- FSL, a comprehensive library of analysis tools for FMRI, MRI and DTI brain imaging data.
- DIAMOND, a c software that is characterizing brain tissue by assessment of the distribution of anisotropic microstructural environments in diffusion-compartment imaging.
- Dmipy, a python library estimating diffusion MRI-based microstructure features, used to fit and recover the parameters of multi-compartment microstructure models

This guide aims to give an introduction to ElikoPy and a brief installation instructions.

INTRODUCTION TO ELIKOPY

ElikoPy is Python library aiming at easing the processing of diffusion imaging for microstructural analysis. ElikoPy expands state of the art pipeline frameworks by providing a complete quality assessment and quality reports for each processed subject, providing a standardized framework to ensure reproducibility and consistency, reducing error propagation and improve sensitivity. It also grants the possibility for clinicians to perform fast preprocessing for a large variety of studies with minimal knowledge.

The ElikoPy library was developed during a Master's thesis.

Note: If you wish to learn more about the library and its validation, we invite you to read our [Master's thesis](#).

1.1 Why use ElikoPy?

Diffusion weighted magnetic resonance imaging (DW-MRI) is a rapidly evolving, non radiating and non invasive technique that allows to capture information on the brain microstructure through the restricted diffusion of water molecules. DW-MRI has seen a growing interest in the recent years motivating the acquisition of large multi-scanner multi-site data sets. The substantial acquisition time of this type of MRI sequence has also encouraged the extensive use of Echo Planar Imaging which suffers from additional artifacts and noise. Several tools have been developed in order to correct those individual problems but they come with the disadvantages of processing only one subject at a time and requiring different softwares making them cumbersome to use. This work presents and evaluates the performances of the ElikoPy library, a complete diffusion MRI processing pipeline that reduces common sources of artifact and captures information on the brain microstructure through multiple microstructural diffusion models. ElikoPy has been designed to deal with large databases and to be robust to different types of acquisitions.

1.2 Features

1. Preprocessing of your dMRI data.
2. Generation of a synthesized b0 for diffusion distortion correction using the based on the [Synb0-DISCO repository](#). This synthesized b0 is usefull for topup if the acquisition was only performed with one phase encoding direction.
3. The library can compute the DTI, Noddi, DIAMOND and the novel Microstucturefingerprinting metric.
4. Complete quality reports to review each step of the processing.
5. Tissue segmentation from T1 images.
6. Ability to run subject and group wise statistics on the dataset.

INSTALLATION

2.1 Manual Installation Steps

You will need a Linux system (CentOS 7 is recommended) to run ElikoPy and all its dependencies natively using a manual installation. Doing a manual installation is not recommended if you have only a limited knowledge in computer science.

2.1.1 Installation of the dependencies

You must first install dependency to your system. Some dependencies are optionnal while other are mandatory.

FSL installation (mandatory)

FSL is a mandatory comprehensive library dependency used among other steps for the preprocessing of diffusion images. FSL is available ready to run on [the official FSL installation page](#).

FreeSurfer installation (optionnal)

FreeSurfer is a software package for the analysis and visualization of structural and functional neuroimaging data from cross-sectional or longitudinal studies. This software is mandatory when correcting from susceptibility distortion using T1 structural images in the preprocessing. To install it, visit the [FreeSurfer Downloads page](#) and pick a package archive suitable to the environment you are in.

ANTs installation (optionnal)

ANTs computes high-dimensional mappings to capture the statistics of brain structure and function. This software is mandatory when correcting from susceptibility distortion using T1 structural images in the preprocessing. ANTs can be compiled from source or installed via pre-built package using their [Github page](#).

C3D installation (optional)

C3D is a command-line tool for converting 3D images between common file formats. The tool also includes a growing list of commands for image manipulation, such as thresholding and resampling. This software is mandatory when correcting from susceptibility distortion using T1 structural images in the preprocessing. A precompiled version of C3D is available on [Sourceforge](#).

Microstructure Fingerprinting installation (recommended) (optional)

Microstructure Fingerprinting estimate the white matter microstructural properties from a dictionary of Monte Carlo diffusion MRI fingerprints. To install it, first download a copy of the [MF repository](#).

```
git clone git@github.com:rensonnetg/microstructure_fingerprinting.git
```

Then, navigate to the folder where this repository was cloned or downloaded (the folder containing the setup.py file) and install the package as follows.

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

DIAMOND installation (optional)

Unfortunately, the DIAMOND code is not publically available. If you do not have it in your possession, you will not be able to use this algorithm. If you have it, simply add the executable to your path.

2.1.2 Installation of ElikoPy

ElikoPy requires Python v3.7+ to run. To install it, first download a copy or clone the [ElikoPy repository](#).

```
git clone git@github.com:Hyedryn/elikopy.git
```

After cloning the repo, you can either firstly install all the python dependencies including optional dependency used to speed up the code.

```
pip install -r requirements.txt --user
```

Or you can install directly the library with only the mandatory dependencies (if you performed the previous step, you still need to perform this step).

```
python3 setup.py install --user
```

Note: When using ElikoPy, do not forget to reference it among all of the used dependencies.

2.2 Container Installation Steps

To ease the installation of ElikoPy, a Singularity container is provided in the [ElikoPy repository](#). To learn more about Singularity, you can visit their [official website](#).

```
git clone https://github.com/Hyedryn/elikopy.git
cd /path/to/repo
sudo singularity build /path/to/elikopy.sif Singularity_elikopy
```

After building the container, ElikoPy can be run using the following command:

```
singularity run -e --contain
-B /path/to/study/directory/:/PROJECTS
-B /tmp:/tmp
-B /path/to/freesurfer/license.txt:/Software/freesurfer/license.txt
-B /path/to/cuda:/usr/local/cuda
--nv
/path/to/elikopy.sif
/path/to/script.py
```

The script.py file contains the Python code that will be executed inside the container. The path to the root directory in your python code must always be “/PROJECTS/” due to the folder binding.

Note: Binding the freesurfer license is optional and is only needed for Synb0-DisCo.

Note: Binding the cuda path is optional and is only needed to speed-up Synb0-DisCo or perform inter slice motion correction with Eddy FSL.

2.3 Using ElikoPy on the CECI Cluster

UCLouvain student who wish to use ElikoPy on the CECI cluster can use the existing installation present in the pilab project directory. First, the following line needs to be added to our *.bash_profile*.

```
source /CECI/proj/pilab/Software/config_elikopy.bash
```

Then, execute the following line of code to install ElikoPy:

```
source /CECI/proj/pilab/Software/install_elikopy.bash
```

If you wish to update your ElikoPy installation, you just need to execute again the preceding line of code.

Authorized user can update the local ElikoPy repository present in the PiLab directory using the following script. The local repository is update using the master branch of the remote Github repository.

```
source /CECI/proj/pilab/Software/update_elikopy.bash
```

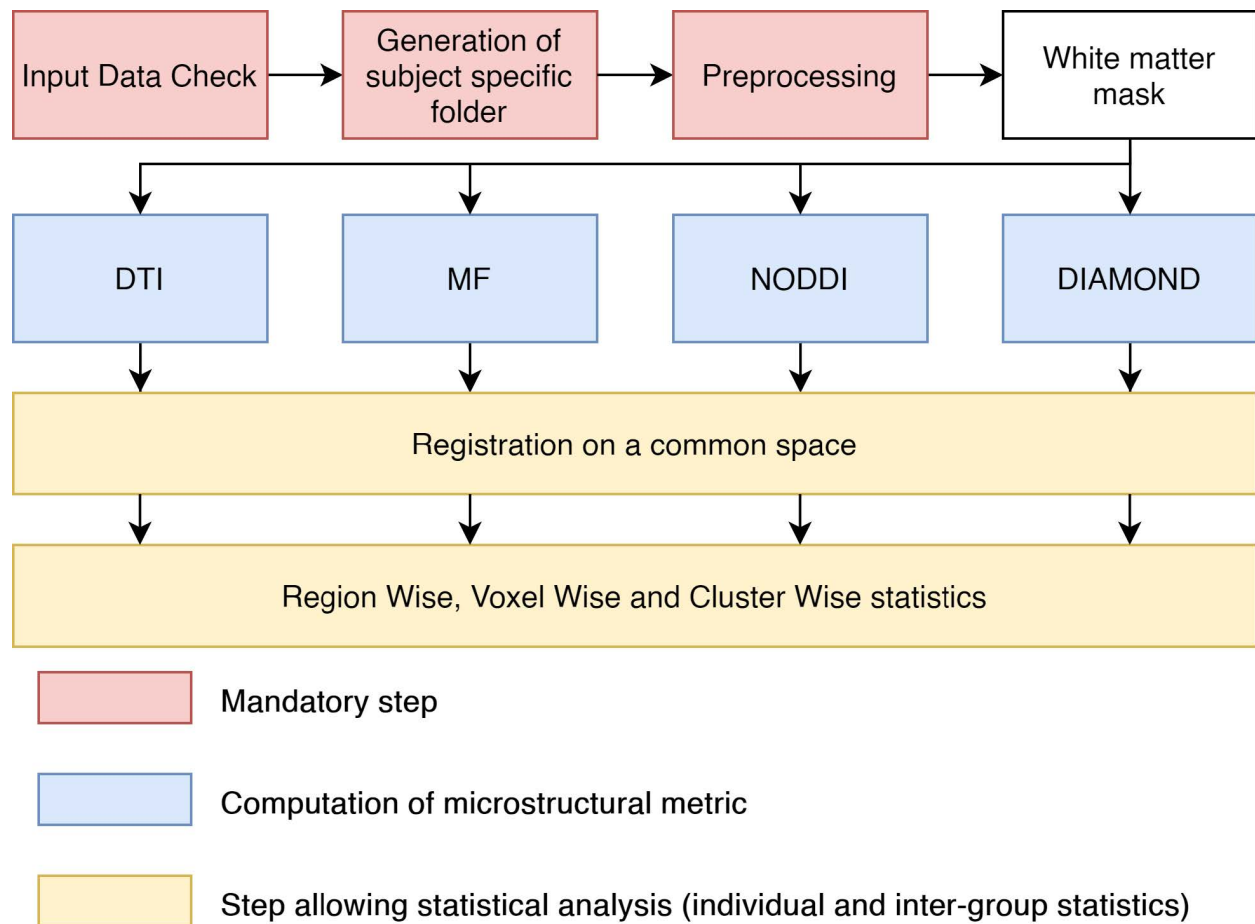
These steps should be sufficient for the *lemaitre3* and *manneback* clusters. When using other clusters, some additional modules may need to be loaded (see the related [CECI documentation](#) for more information). We also strongly recommend you to familiarize yourself with slurm job when using ElikoPy on the CECI cluster.

PROJECT AS AN EASY WAY TO MANAGE A STUDY

3.1 Main available processing steps

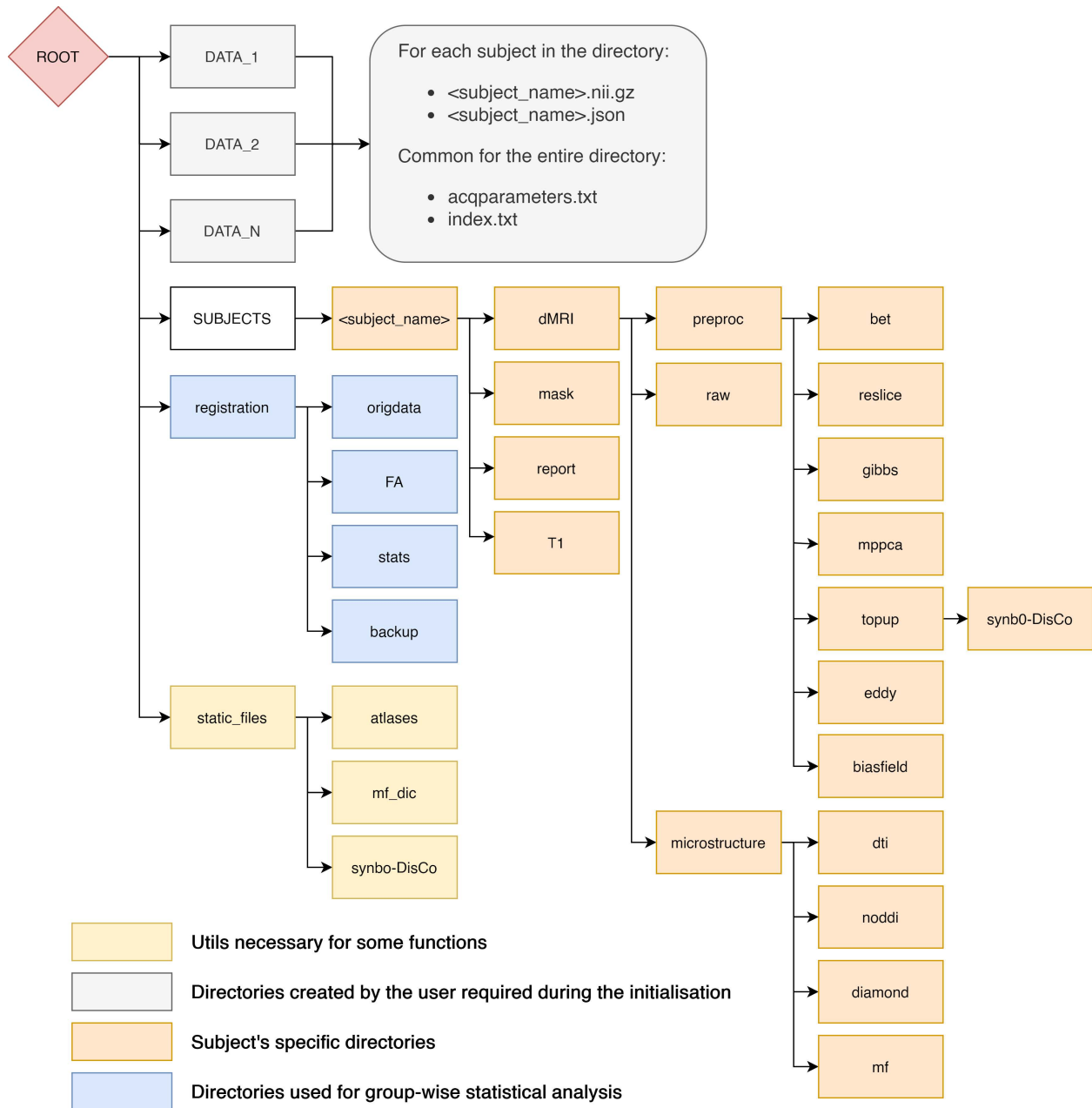
Using a rich set of functions, data processed with ElikoPy are firstly audited to ensure that the image dimension match the dimension of the b-value and b-vector files along the acquisitions parameters and index files. If successful, a dedicated storage folder for each subject is generated. Afterward, preprocessing can be applied to correct and enhanced the raw image. This includes, brain extraction, reslicing, Principal Component Analysis (PCA) based denoising using the Marchenko-Pastur distribution, suppression in image space of Gibbs ringing artifacts, estimation and correction of the susceptibility induced field, modeling and correction of subject movements and finally, bias field correction. After the preprocessing, a white matter mask registered on diffusion data can be obtained using a T1 image or computed directly from diffusion data by segmenting an Anisotropic Power (AP) map.

As seen in the figure below, four distinct algorithms for the estimation of the microstructure are available. Subsequently, the pipeline adds the possibility to register the computed metrics in a common space given by either one of the study subjects or by an atlas. Finally, from there, ElikoPy can output statistical results for population studies by aligning metrics from multiple subjects into a common space and performing region wise, voxel wise and cluster wise statistics.



3.2 Folder structure

The Elikopy toolbox follows a specific folder structure that prevents ambiguity and data losses when dealing with a large amount of subjects. The figure below illustrates the folder tree in ElikoPy.



The first type of folder present at the root of the Elikopy project are **data** folders. These folders contain all raw subject files belonging to the same class along with the associated acquisition parameters and index files. Classes are used to separate subjects that have different acquisition parameters or subjects that need to be separated from others groups of subjects. The pipeline does not have a limitation on the number of classes.

The **subjects** folder contains a subfolder for each valid subject presents in **data** folders. Along these subfolders, three json files are present. The `subj_error.json` file contains the list of subjects with invalid raw data, The `subj_list.json` file contains the list of valid subjects and the `subj_type.json` is a dictionary that maps each subjects subfolder to its data class.

Each subdirectory of the **subjects** folder contains the output of every preprocessing and processing function executed on the patient associated with the subdirectory. The output consists of NIfTI files, log files and some others files related to the specific functions.

The **registration** folder contains diffusion metrics registered to a common space, group wise statistics and voxel wise

statistics for each registered metric.

Finally, the **static_files** folder contains files mandatory for some processing steps of the library such as MF dictionary and Synb0-DisCo atlases.

TYPICAL USAGE FOR PROCESSING A STUDY

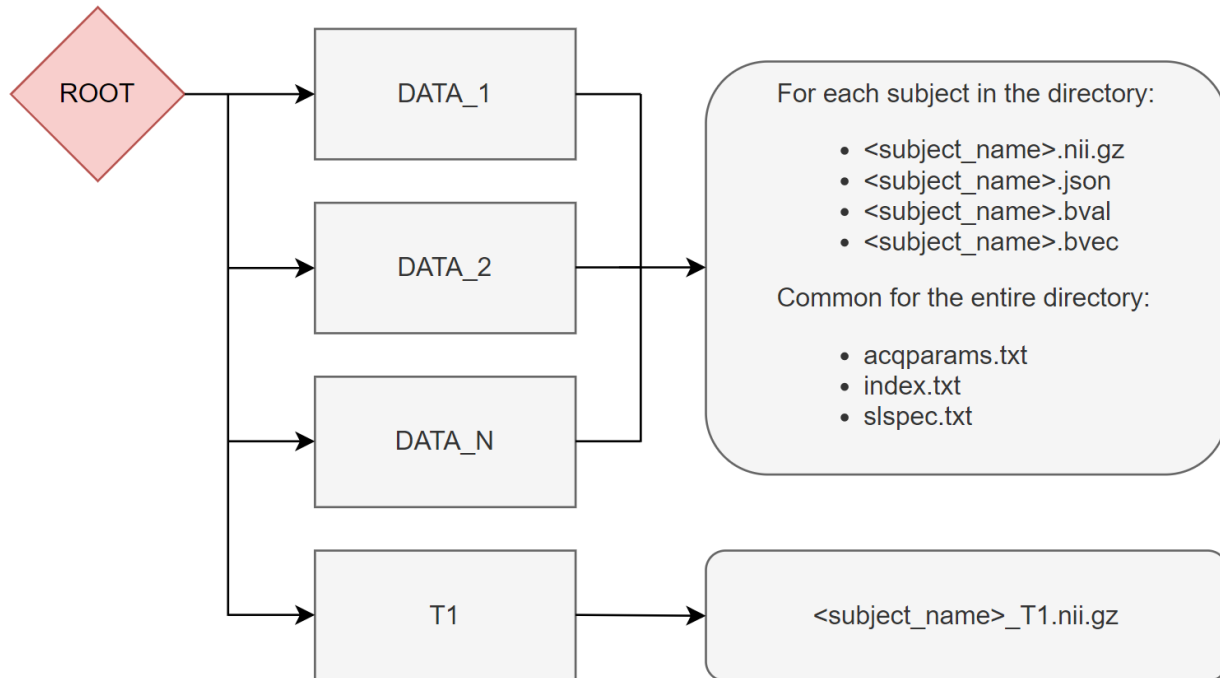
On this page is presented a basic usage of the ElikoPy library. More information on all these functions are available in the detailed guide.

4.1 Header and initialisation

The first step to enable ElikoPy is to import it and initialise the ElikoPy object “*study*” specific to the current study. The only required argument for the constructor is the path to the root directory of the project.

```
1 import elikopy
2 import elikopy.utils
3
4 f_path="/PROJECTS/"
5 dic_path="/PROJECTS/static_files/mf_dic/fixed_rad_dist.mat"
6
7 study = elikopy.core.Elikopy(f_path)
8 study.patient_list()
```

The root directory must have the following structure during the initialisation



The T1 structural images as well as the acqparams, index and slspec files are optional. However, if they are not available, some processing steps might be not available (this is usually specified by a note in the documentation). These files can be generated as explained in the following links:

- acqparams.txt and index.txt : [Eddy FSL acqp](#)
- slspec.txt : [Eddy FSL slspec](#)

4.2 Preprocessing

The following code block show how to preprocess the dMRI data. By default only the brain extraction is enabled in the preprocessing but we recommend you to enable more preprocessing as described in the detailed guide (see [Preprocessing of diffusion images](#)).

```
8 study.preproc()
```

4.3 whitematter mask

The following code block computes a white matter mask for each subject from its T1 structural image (if available). If the T1 is not available, the mask is computed using the anisotropic power map generated from the diffusion data.

```
9 study.white_mask()
```

4.4 Microstructural metrics computation

The following code block computes microstructural metrics from the four microstructural model available in ElikoPy.

```

10 study.dti()
11 study.noddi()
12 study.diamond()
13 study.fingerprinting()

```

4.5 Statistical Analysis

In the following code block, fractional anisotropy (FA) from DTI along other additional metrics are registered into a common space. The registration is computed using the FA and the mathematical transformation is applied to other metrics.

Afterwards, the `randomise_all` function performs group wise statistic for the defined metrics along extraction of individual region wise value for each subject into csv files.

```

14 grp1=[1]
15 grp2=[2]
16
17
18
19 study.regall_FA(grp1=grp1,grp2=grp2)
20
21 additional_metrics={'_noddi_odi':'noddi','_mf_fvf_tot':'mf','_diamond_kappa':'diamond'}
22 study.regall(grp1=grp1,grp2=grp2, metrics_dic=additional_metrics)
23
24 metrics={'dti':'FA','_noddi_odi':'noddi','_mf_fvf_tot':'mf','_diamond_kappa':'diamond'}
25 study.randomise_all(metrics_dic=metrics)

```

4.6 Data Exportation

The export function is used to “revert” the folder structure, instead of using a subject specific folder tree, data are exported into a metric specific folder tree. In this example, only metrics computed from the dti model are exported.

```

22 study.export(raw=False, preprocessing=False, dti=True,
23             noddi=False, diamond=False, mf=False, wm_mask=False, report=True)

```

Note: If you wish to learn more about the library and its validation, we recommend you to read the detailed guide and play around with the library.

4.7 Other parameters commonly available

The ElikoPy library has been made compatible with the slurm scheduler commonly present on HPC clusters. The use of the slurm scheduler can be controlled with the **slurm** parameters.

Associated options are:

- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – If not None, Topup will use additional parameters based on the supplied config file located at <topupConfig>. default=None Email adress to send notification if a task fails. default=None
- **slurm_timeout** - Replace the default slurm timeout used in the ElikoPy function by a custom timeout.
- **slurm_mem** - Replace the default amount of ram allocated to the slurm task by a custom amount of ram.
- **cpus** – Replace the default number of slurm cpus by a custom number of cpus.

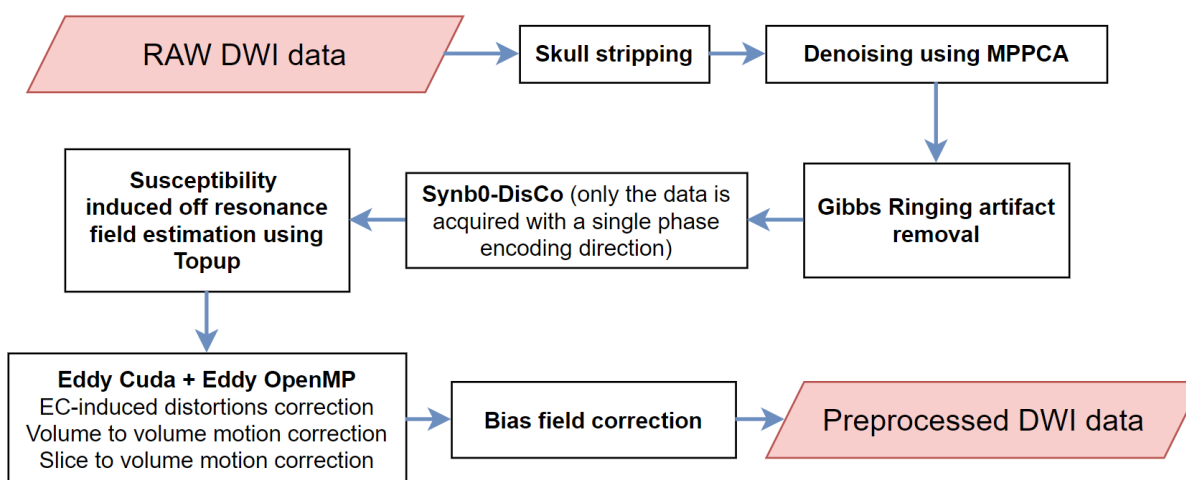
The slurm option and slurm_email option can be globally define during the initialisation of the study object.

When processing a study, the processing for some subjects could fail for various reasons. The ElikoPy library provides two parameters destined to limit the amount of processing necessary to recover from these failures.

- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **starting_state** – Manually set which step of the function to start from. default=None

PREPROCESSING OF DIFFUSION IMAGES

The preprocessing stage aims at correcting the non idealities affecting the diffusion data before computing the diffusion metrics. With the exception of skull stripping, all processing steps are optional and can be applied at the user discretion.



To preprocess the dMRI data, the following line of code is used. However, this results in the default preprocessing that encompass only the skull stripping step. To perform more advanced preprocessing, we need to dive into the arguments of the preproc function.

```
study.preproc()
```

The arguments of the preproc function are given in the API : [LINK](#). In this page, only the main arguments are explained in order to grasp the key aspects of preprocessing using ElikoPy.

5.1 Reslice

5.1.1 Description

If the raw data is not in its 'native' resolution, a reslicing process might be required. Usually, the MRI scanner performs automatic interpolation on the data in order to beautify the data since clinicians usually have a preference for high resolution images. However, the intrinsic resolution is not augmented by this interpolation. While somewhat useful to clinicians, the interpolation is usually not desirable for research. Using it means more computation time and uncorrelated noise becoming correlated which reduces the performances of MPPCA denoising algorithms. Moreover, interpolation is not desirable when performing Gibbs ringing correction. Reslicing is therefore a way to mitigate the effect of mandatory interpolation during the acquisition.

5.1.2 Related parameters

The reslicing step during the preprocessing can be activated using the `reslice` argument.

- **reslice** - If true, data will be resliced with a new voxel resolution of $2 \times 2 \times 2$. default=False
- **reslice_addSlice** - If true, an additional empty slice will be added to each volume (might be useful for motion correction if one slice is dropped during the acquisition and the user still wants to perform easily the slice-to-volume motion correction). default=False

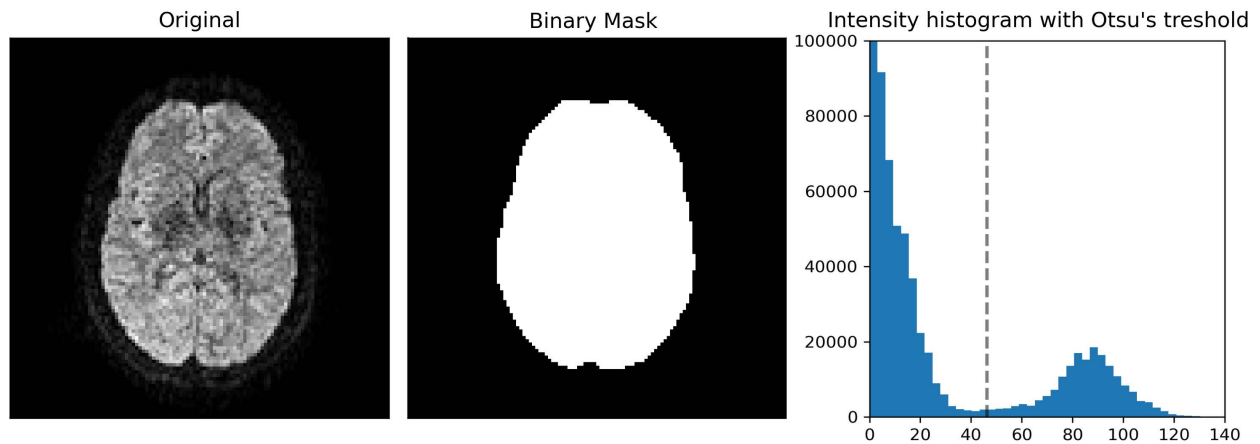
```
study.preproc(reslice=True, reslice_addSlice=False)
```

5.2 Brain Extraction

5.2.1 Description

The brain is extracted from the skull and other tissues surrounding the brain to increase the processing efficiency of subsequent steps and it is generally required before using other image processing algorithms. At the end of the preprocessing, a final brain mask readjusted in regard of all the applied preprocessing steps is also provided as output.

The mask is computed using `median_otsu` from DiPy.



5.2.2 Related parameters

The brain extraction is the only mandatory step and cannot be disabled. However, it is possible to change the parameters of the method

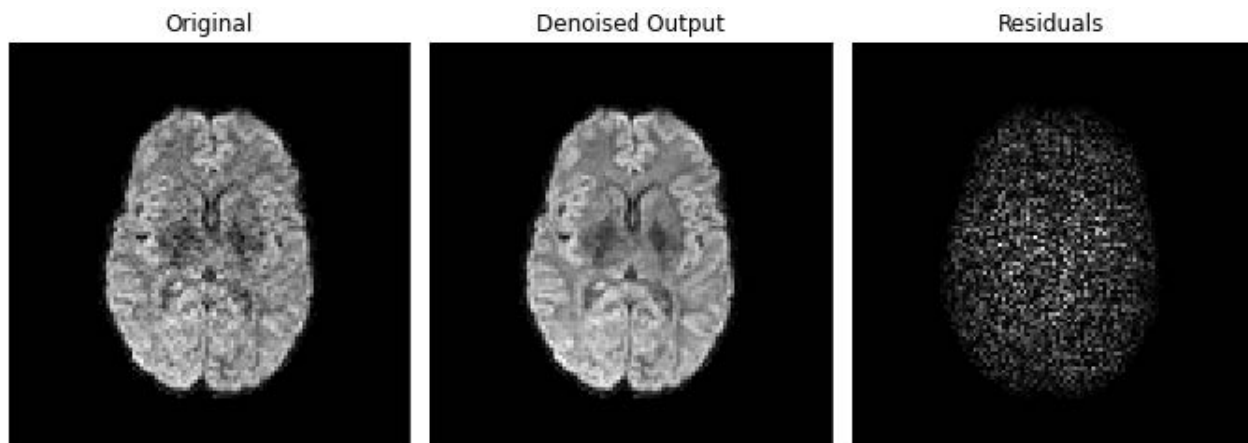
- **bet_median_radius** - Radius (in voxels) of the applied median filter during brain extraction. default=2
- **bet_numpass** - Number of pass of the median filter during brain extraction. default=1
- **bet_dilate** - Number of iterations for binary dilation during brain extraction. default=2

```
study.preproc(bet_median_radius=2, bet_numpass=2, bet_dilate=2)
```

5.3 MPPCA Denoising

5.3.1 Description

To reduce Rician noise typically found in MR images, the input images are denoised using the Marchenko-Pastur PCA technique as implemented in DiPy. Since the noise in diffusion data is spatially dependent in the case of multichannel receive coils, Principal component analysis of Marchenko-Pastur (MPPCA) noise-only distribution provides an accurate and fast method of noise evaluation and reduction. This methods has been chosen since it is a fast denoising algorithm that does not blur the image or create artifact.



5.3.2 Related parameters

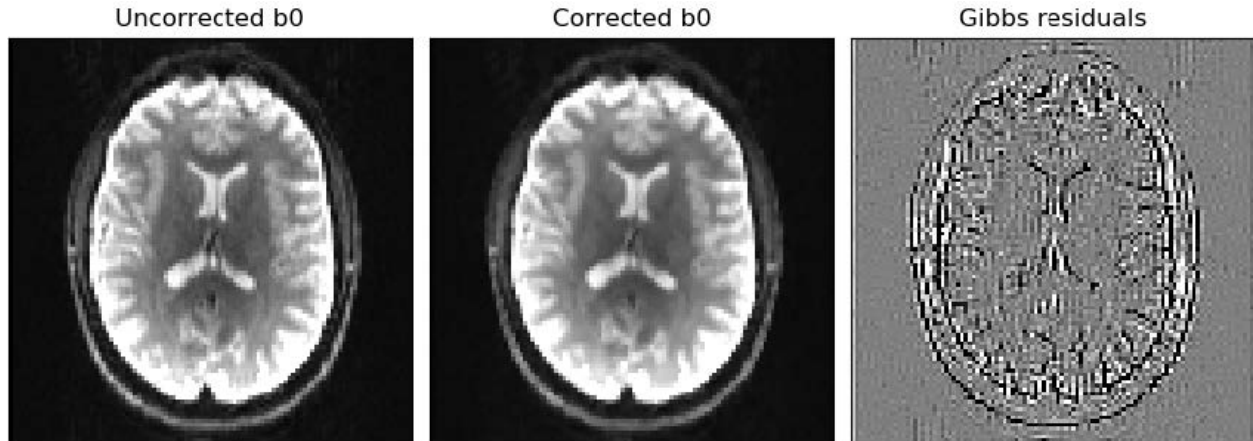
The denoising step during the preprocessing can be activated using the denoising argument.

```
study.preproc(denoising=True)
```

5.4 Gibbs Ringing Correction

5.4.1 Description

In general, in the context of diffusion-weighted imaging, derived diffusion-based estimates are affected by Gibbs oscillations. To correct for this, `gibbs_removal` from DiPy is used. This algorithm models the truncation of k-space as a convolution with a sinc-function in the image space. The severity of ringing artifacts thus depends on how the sampling of the sinc function occurs. The `gibbs_removal` function reinterpolate the image based on local, subvoxel-shifts to sample the ringing pattern at the zero-crossings of the oscillating sinc-function.



5.4.2 Related parameters

The Gibbs removal can be enabled using the `gibbs` argument.

```
study.preproc(gibbs=True)
```

Unless the data suffer heavily from Gibbs ringing artifacts, we do not advise to use the gibbs ringing removal step as it might blurr out small microstructural features.

5.5 Susceptibility field estimation

5.5.1 Description

Susceptibility distortions are created by differences in magnetic susceptibility near junctions of tissues. The susceptibility off resonance field is estimated using Topup from FSL. To do so, Topup needs data acquired with multiple phase encoding directions (at least 2). If only a single phase encoding direction is available, ElikoPy uses instead a generated synthetic volume based on a T1 structural image using Synb0-DisCo. This step only allows to **estimate** the susceptibility distortions, they are corrected at the same time as the eddy current distortions in the Eddy step below.

5.5.2 Related parameters

The susceptibility field estimation can be enabled using the `topup` argument.

- **topup** - true, Topup will estimate the susceptibility induced distortions. These distortions are corrected at the same time as EC-induced distortions if `eddy=True`. In the absence of images acquired with a reverse phase encoding direction, a T1 structural image is required. `default=False`
- **topupConfig** – If not None, Topup will use additionnal parameters based on the supplied config file located at `<topupConfig>`. `default=None`
- **forceSynb0DisCo** - If true, Topup will always estimate the susceptibility field using the T1 structural image. `default=False`

```
study.preproc(topup=True)
```

Note: If Topup is used, ElikoPy needs the acqparam and index files when generating the patient list : [LINK](#) (page getting started)

Note: If topup is enabled for data with a single phase encoding direction, a T1 structural image has to be provided when generating the patient list : [LINK](#) (page getting started)

5.6 Eddy and motion correction

5.6.1 Description

Motion, susceptibility and Eddy current induced distortions are artifacts with different origins but a similar effect i.e the displacement and deformation of the brain. They can therefore be jointly corrected. This is achieved using FSL Eddy. The susceptibility distortions are only corrected if they have been estimated during the topup step. By default only the inter-volume (volume-to-volume) motion is corrected but it is also possible to correct for intra-volume (slice-to-volume) motion.

5.6.2 Related parameters

The motion and distortion correction can be activated using the eddy argument. The number of iteration for the motion correction algorithm can also be changed.

```
study.preproc(eddy=True, niter=5)
```

In cases with large motion, inter-volume motion correction might not be sufficient and intra-volume correction is required. This option can be enabled using the s2v argument. The s2v input is a list of 4 parameters : [mporder,s2v_niter,s2v_lambda,s2v_interp]. The slice-to-volume motion correction is performed if mporder>0. These parameters are explained in depth in the FSL documentation ([LINK](#)). If N describes the number of excitations in a volume, setting mporder to N/4 while letting the other 3 parameters to their default values should provide good results in most situations. The slice-to-volume motion correction is only possible with cuda enabled.

Using the framework of Eddy FSL, it is also possible to replace outlier slices. This is done with the olrep argument which is a list of 4 parameters : [repol,ol_nstd,ol_nvox,ol_type]. The outlier replacement is performed if repol==True. These parameters are explained in depth in the FSL documentation.

```
study.preproc(eddy=True, niter=5, s2v=[6,5,1,'trilinear'], cuda=True, cuda_name='eddy_
↪cuda10.1', olrep=[True, 4, 250, 'sw'])
```

Note: If Eddy FSL is used, ElikoPy needs the acqparam and index files when generating the patient list : [LINK](#) (page getting started)

Note: If slice-to-volume motion correction is enabled, ElikoPy needs the slspec file when generating the patient list : [LINK](#) (page getting started)

5.7 Bias Field Correction

5.7.1 Description

Variability of the signal in tissues of the same type can affect microstructural metrics computation and brain segmentation algorithms. This can be corrected using the N4 Bias Field Correction algorithm.

5.7.2 Related parameters

The bias field correction can be activated using the `biasfield` argument. It is also possible to modify the parameters of the correction method.

- **`biasfield_bsplineFitting`** - Define the initial mesh resolution in mm and the bspline order of the biasfield correction tool.
- **`biasfield_convergence`** - Define the maximum number of iteration and the convergences threshold of the biasfield correction tool.

```
study.preproc(biasfield=True, biasfield_bsplineFitting=[100,3], biasfield_
↳ convergence=[1000,0.001])
```

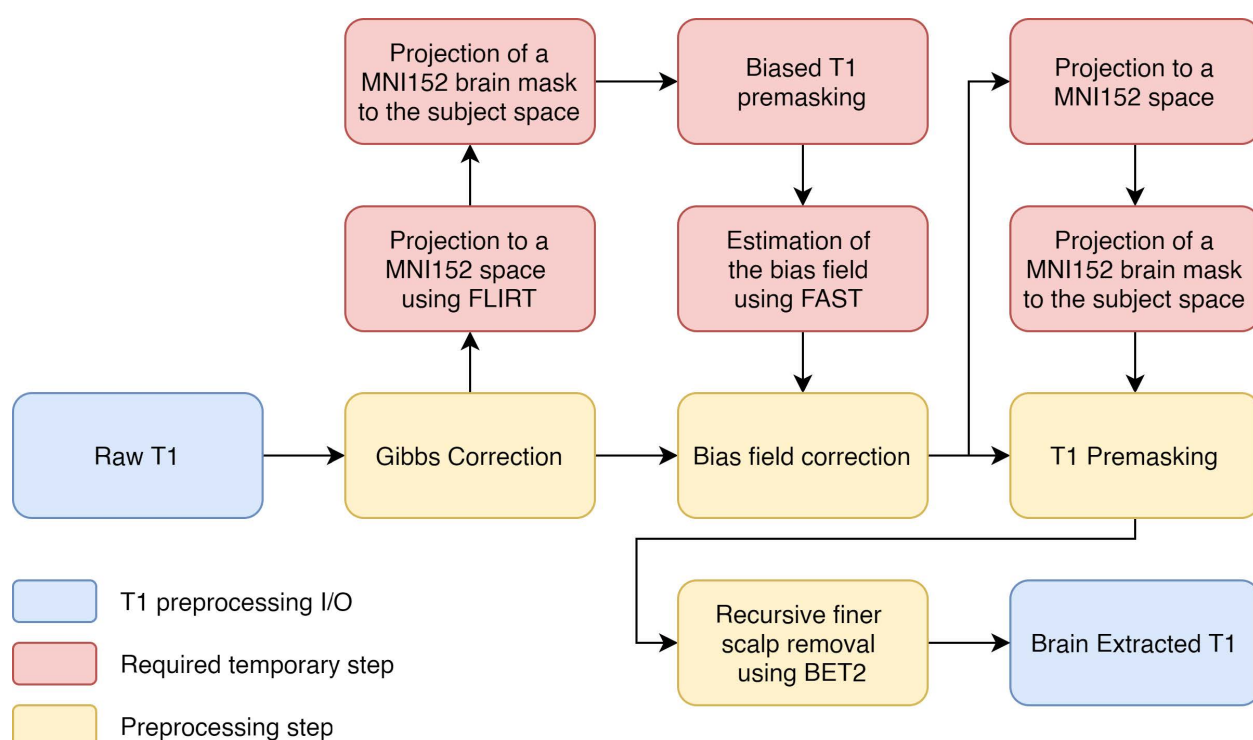
5.8 Report

By default, the `preproc` function outputs a quality report that contains quality control features for the processing. This can be disabled if needed.

```
study.preproc(report=False)
```

T1 PREPROCESSING

Providing a white matter mask is a useful step to accelerate microstructural features computation and more easily do tractography. The `white_mask` function of the ElikoPy library has been elaborated to perform this important step.



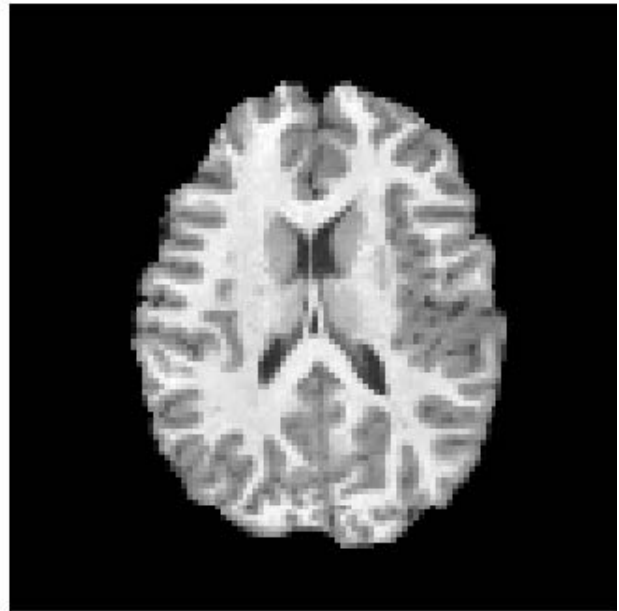
On the one hand, when a T1 image is available, a white matter mask can be computed from this data. Therefore, the T1 image is first preprocessed then segmented. Finally the segmented white matter mask is projected into the space of the preprocessed diffusion image.

On the other hand, when no T1 images are available, the white matter mask is directly computed from a segmentation of the diffusion data using Anisotropic Power (AP) map. In this case, no registrations are necessary.

Anisotropic Power Map (pseudo T1)



Unbiased Gibbs Corrected T1



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EXAMPLES

CONTRIBUTING

ElikoPy is an open source project, meaning we have the challenge of limited resources. We are grateful for any support that you can offer. Helping other users, raising issues, helping write documentation, or contributing code are all ways to help!

11.1 Raise an Issue

For general bugs/issues, you can open an issue [at the GitHub repo](#).

11.2 Write Documentation

We (like almost all open source software providers) have a documentation dilemma. . . We tend to focus on the code features and functionality before working on documentation. And there is very good reason for this: we want to share the love so nobody feels left out!

You can contribute to the documentation by [raising an issue to suggest an improvement](#) or by sending a [pull request](#) on [our repository](#).

11.3 Contribute to the code

We use the traditional [GitHub Flow](#) to develop. This means that you fork the main repo, create a new branch to make changes, and submit a pull request (PR) to the master branch.

11.3.1 Step 1. Fork the repo

To contribute to ElikoPy, you should obtain a GitHub account and fork the [ElikoPy](#) repository. Once forked, clone your fork of the repo to your computer. (Obviously, you should replace `your-username` with your GitHub username.)

```
$ git clone https://github.com/your-username/elikopy.git && \  
  cd elicopy/
```

11.3.2 Step 2. Checkout a new branch

Branches are a way of isolating your features from the main branch. Given that we've just cloned the repo, we will probably want to make a new branch from master in which to work on our new feature. Lets call that branch `new-feature`:

```
$ git checkout master && \  
    git checkout -b new-feature
```

Note: You can always check which branch you are in by running `git branch`.

11.3.3 Step 3. Make your changes

On your new branch, go nuts! Make changes, test them, and when you are happy commit the changes to the branch:

```
$ git add file-changed1 file-changed2...  
  
$ git commit -m "what changed?"
```

This commit message is important - it should describe exactly the changes that you have made. Good commit messages read like so:

```
$ git commit -m "changed function preproc in core.py to output new mask to fix #2"  
  
$ git commit -m "updated docs about MF to close #10"
```

The tags `close #10` and `fix #2` are referencing issues that are posted on the upstream repo where you will direct your pull request. When your PR is merged into the master branch, these messages will automatically close the issues, and further, they will link your commits directly to the issues they intend to fix. This will help future maintainers understand your contribution, or (hopefully not) revert the code back to a previous version if necessary.

11.3.4 Step 4. Push your branch to your fork

When you are done with your commits, you should push your branch to your fork (and you can also continuously push commits here as you work):

```
$ git push origin new-feature
```

Note that you should always check the status of your branches to see what has been pushed (or not):

```
$ git status
```

11.3.5 Step 5. Submit a Pull Request

Once you have pushed your branch, then you can go to your fork (in the web GUI on GitHub) and [submit a Pull Request](#). Regardless of the name of your branch, your PR should be submitted to the ElikoPy master branch. Submitting your PR will open a conversation thread for the maintainers of ElikoPy to discuss your contribution. At this time, the continuous integration that is linked with the code base will also be executed. If there is an issue, or if the maintainers suggest changes, you can continue to push commits to your branch and they will update the Pull Request.

11.3.6 Step 6. Keep your branch in sync

Cloning the repo will create an exact copy of the ElikoPy repository at that moment. As you work, your branch may become out of date as others merge changes into the upstream master. In the event that you need to update a branch, you will need to follow the next steps:

```
$ git remote add upstream https://github.com/Hyedryn/elikopy.git && # to add a new ↵
↵remote named "upstream" \
    git checkout master && # or another branch to be updated \
    git pull upstream master && \
    git push origin master && # to update your fork \
    git checkout new-feature && \
    git merge master
```


ELIKOPY PACKAGE

12.1 Submodules

12.2 `elikopy.core` module

Elikopy @author: qdessain, msimon

class `elikopy.core.Elikopy`(*folder_path*, *cuda=False*, *slurm=False*, *slurm_email='example@example.com'*)
Bases: `object`

Main class containing all the necessary function to process and preprocess a specific study.

diamond(*folder_path=None*, *patient_list_m=None*, *reportOnly=False*, *slurm=None*, *slurm_email=None*,
slurm_timeout=None, *cpus=None*, *slurm_mem=None*)

Computes the DIAMOND metrics for each subject. The outputs are available in the directories
<folder_path>/subjects/<subjects_ID>/dMRI/microstructure/diamond/.

example : `study.diamond()`

Parameters

- **folder_path** – the path to the root directory. default=`study_folder`
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : [`'patientID1'`, `'patientID2'`, `'patientID3'`]. default=`None`
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=`value_during_init`
- **slurm_email** – Email adress to send notification if a task fails. default=`None`
- **slurm_timeout** – Replace the default slurm timeout of 14h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 4 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (6096MO by cpu) by a custom amount of ram.

dti(*folder_path=None*, *patient_list_m=None*, *slurm=None*, *slurm_email=None*, *slurm_timeout=None*,
slurm_cpus=None, *slurm_mem=None*)

Computes the DTI metrics for each subject using Weighted Least-Squares. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/dMRI/dti/.

example : `study.dti()`

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 1h by a custom timeout.
- **slurm_cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slurm, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

export(*folder_path=None, raw=False, preprocessing=False, dti=False, noddi=False, diamond=False, mf=False, wm_mask=False, report=False, preprocessed_first_b0=False, patient_list_m=None, tractography=False*)

Allows to obtain in a single Export folder the outputs of specific processing steps for all subjects.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **raw** – If true, copy the raw data of each subject in the Export folder. default=FALSE
- **preprocessing** – If true, copy the preprocessed data of each subject in the Export folder. default=FALSE
- **dti** – If true, copy the DTI outputs of each subject in the Export folder. default=FALSE
- **noddi** – If true, copy the NODDI outputs of each subject in the Export folder. default=FALSE
- **diamond** – If true, copy the DIAMOND outputs of each subject in the Export folder. default=FALSE
- **mf** – If true, copy the MF outputs of each subject in the Export folder. default=FALSE
- **wm_mask** – If true, copy the white matter mask of each subject in the Export folder. default=FALSE
- **report** – If true, copy the quality control reports of each subject in the Export folder. default=FALSE
- **tractography** – If true, copy the tractography outputs of each subject in the Export folder. default=FALSE

fingerprinting(*dictionary_path=None, folder_path=None, CSD_bvalue=None, slurm=None, patient_list_m=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Computes the Microstructure Fingerprinting metrics for each subject. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/dMRI/microstructure/mf/.

example : `study.fingerprinting(dictionary_path='my_dictionary')`

Parameters

- **folder_path** – the path to the root directory. default=`study_folder`
- **dictionary_path** – Path to the dictionary of fingerprints (mandatory).
- **CSD_bvalue** – If the DIAMOND outputs are not available, the fascicles directions are estimated using a CSD with the images at the b-values specified in this argument. default=`None`
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : `['patientID1','patientID2','patientID3']`. default=`None`
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=`value_during_init`
- **slurm_email** – Email adress to send notification if a task fails. default=`None`
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **slurm_cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

noddi(*folder_path=None, patient_list_m=None, force_brain_mask=False, slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Computes the NODDI metrics for each subject. The outputs are available in the directories `<folder_path>/subjects/<subjects_ID>/dMRI/microstructure/noddi/`.

example : `study.noddi()`

Parameters

- **folder_path** – the path to the root directory. default=`study_folder`
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : `['patientID1','patientID2','patientID3']`. default=`None`
- **force_brain_mask** – Force the use of a brain mask even if a whitematter mask exist. default=`False`
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=`value_during_init`
- **slurm_email** – Email adress to send notification if a task fails. default=`None`
- **slurm_timeout** – Replace the default slurm timeout of 10h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

noddi_amico(*folder_path=None, patient_list_m=None, force_brain_mask=False, slurm=None, slurm_email=None, slurm_timeout=None, slurm_cpus=None, slurm_mem=None*)

Computes the NODDI amico metrics for each subject. The outputs are available in the directories `<folder_path>/subjects/<subjects_ID>/dMRI/microstructure/noddi/`.

example : `study.noddi_amico()`

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **force_brain_mask** – Force the use of a brain mask even if a whitematter mask exist. default=False
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 10h by a custom timeout.
- **slurm_cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slurm, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

noddi_fix_icvf_thresholding(*folder_path=None, patient_list_m=None, fintra_threshold=0.99, fbundle_threshold=0.05, use_brain_mask=False, use_wm_mask=False*)

A function to quickly change the treshold value applied on the icvf metric of noddi without the needs of executing again the full noddi core function.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **fintra_threshold** – Threshold applied on the fintra. default=0.99
- **fbundle_threshold** – Threshold applied on the fbundle. default=0.05
- **use_brain_mask** – Set to 0 values outside the brain mask. default=False
- **use_wm_mask** – Set to 0 values outside the white matter mask. default=False

patient_list(*folder_path=None, bids_path=None, reverseEncoding=True*)

From the root folder containing data_1, data_2, ... data_n folders with nifti files (and their corresponding bvals and bvecs), the Elikopy folder structure is created in a directory named 'subjects' inside folder_path. This step is mandatory. The validity of all the nifti present in the root folder is verified. If some nifti do not possess an associated bval and bvec file, they are discarded and the user is notified in a summary file named subj_error.json generated in the out sub-directory. All valid patients are stored in a file named patient_list.json. In addition to the nifti + bval + bvec, the data_n folders can also contain the json files (with the patient informations) as well as the acquparam, index and slspec files (used during the preprocessing). If these files are missing a warning is raised. In addition to the DW images, T1 structural images can be provided in a directory called 'T1' in the root folder.

example : study.patient_list()

Parameters

- **folder_path** – Path to the root folder of the study. default = study_folder
- **bids_path** – Path to the optional folder containing subjects' data in the BIDS format.
- **reverseEncoding** – Append reverse encoding direction to the DW-MRI data if available. default = True

patientlist_wrapper(*function, func_args, folder_path=None, patient_list_m=None, filename=None, function_name=None, slurm=False, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

A wrapper function that apply a function given as an argument to every subject of the study. The wrapped function must takes two arguments as input, the patient_name and the path to the root of the study.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **function** – The pointer to the function (only without slurm !)
- **func_args** – Additional arguments to pass to the wrapped function (only without slurm !)
- **filename** – The name of the file containing the wrapped function (only with slurm !)
- **function_name** – The name of the wrapped function (only with slurm !)
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slurm, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

preproc(*folder_path=None, reslice=False, reslice_addSlice=False, denoising=False, gibbs=False, topup=False, topupConfig=None, forceSynb0DisCo=False, useGPUSynb0DisCo=False, eddy=False, biasfield=False, biasfield_bsplineFitting=[100, 3], biasfield_convergence=[1000, 0.001], patient_list_m=None, starting_state=None, bet_median_radius=2, bet_numpass=1, bet_dilate=2, cuda=None, cuda_name='eddy_cuda10.1', s2v=[0, 5, 1, 'trilinear'], olrep=[False, 4, 250, 'sw'], slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None, qc_reg=True, niter=5, slspec_gc_path=None, report=True*)

Performs data preprocessing. By default only the brain extraction is enabled. Optional preprocessing steps include : reslicing, denoising, gibbs ringing correction, susceptibility field estimation, EC-induced distortions and motion correction, bias field correction. The results are stored in the preprocessing subfolder of each study subject <folder_path>/subjects/<subjects_ID>/dMRI/preproc.

example : study.preproc(denoising=True, topup=True, eddy=True, biasfield=True)

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **reslice** – If true, data will be resliced with a new voxel resolution of 2*2*2. default=False
- **reslice_addSlice** – If true, an additional empty slice will be added to each volume (might be useful for motion correction if one slice is dropped during the acquisition and the user still wants to perform easily the slice-to-volume motion correction). default=False
- **denoising** – If true, MPPCA-denoising is performed on the data. default=False
- **gibbs** – If true, Gibbs ringing correction is performed. We do not advise to use this correction unless the data suffers from a lot of Gibbs ringing artifacts. default=False

- **topup** – If true, Topup will estimate the susceptibility induced distortions. These distortions are corrected at the same time as EC-induced distortions if eddy=True. In the absence of images acquired with a reverse phase encoding direction, a T1 structural image is required. default=False
- **topupConfig** – If not None, Topup will use additional parameters based on the supplied config file located at <topupConfig>. default=None
- **forceSynb0DisCo** – If true, Topup will always estimate the susceptibility field using the T1 structural image. default=False
- **eddy** – If true, Eddy corrects the EC-induced (+ susceptibility, if estimated) distortions and the motion. If these corrections are performed the acquparam and index files are required (see documentation). To perform the slice-to-volume motion correction the slspec file is also needed. default=False
- **biasfield** – If true, low frequency intensity non-uniformity present in MRI image data known as a bias or gain field will be corrected. default=False
- **biasfield_bsplineFitting** – Define the initial mesh resolution in mm and the bspline order of the biasfield correction tool. default=[100,3]
- **biasfield_convergence** – Define the maximum number of iteration and the convergences threshold of the biasfield correction tool. default=[1000,0.001]
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **starting_state** – Manually set which step of the preprocessing to execute first. Could either be None, denoising, gibbs, topup, eddy, biasfield, report or post_report. default=None
- **bet_median_radius** – Radius (in voxels) of the applied median filter during brain extraction. default=2
- **bet_numpass** – Number of pass of the median filter during brain extraction. default=1
- **bet_dilate** – Number of iterations for binary dilation during brain extraction. default=2
- **cuda** – If true, eddy will run on cuda with the command name specified in cuda_name. default=False
- **cuda_name** – name of the eddy command to run when cuda==True. default="eddy_cuda10.1"
- **s2v** – list of parameters of Eddy for slice-to-volume motion correction (see Eddy FSL documentation): [mporder,s2v_niter,s2v_lambda,s2v_interp]. The slice-to-volume motion correction is performed if mporder>0, cuda is used and a slspec file is provided during the patient_list command. default=[0,5,1,'trilinear']
- **olrep** – list of parameters of Eddy for outlier replacement (see Eddy FSL documentation): [repol,ol_nstd,ol_nvox,ol_type]. The outlier replacement is performed if repol==True. default=[False, 4, 250, 'sw']
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout by a custom timeout.
- **cpus** – Replace the default number of slurm cpus by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.

- **slurm_mem** – Replace the default amount of ram allocated to the slurm task by a custom amount of ram.
- **qc_reg** – If true, the motion registration step of the quality control will be performed. We do not advise to use this argument as it increases the computation time. default=False
- **niter** – Define the number of iterations for eddy volume-to-volume. default=5
- **slspec_gc_path** – Path to the folder containing volume specific slice-specification for eddy. If not None, eddy motion correction with gradient cycling will be performed.
- **report** – If False, no quality report will be generated. default=True

randomise_all(*folder_path=None, randomise_numberofpermutation=5000, skeletonised=True, metrics_dic={'FA': 'dti', '_diamond_kappa': 'diamond', '_mf_fvf_tot': 'mf', '_noddi_odi': 'noddi'}, regionWiseMean=True, slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Performs tract base spatial statistics (TBSS) between the data in grp1 and grp2 (groups are specified during the call to regall_FA) for each diffusion metric specified in the argument metrics_dic. The mean value of the diffusion metrics across atlases regions can also be reported in CSV files using the regionWiseMean flag. The used atlases are : the Harvard-Oxford cortical and subcortical structural atlases, the JHU DTI-based white-matter atlases and the MNI structural atlas It is mandatory to have performed regall_FA prior to randomise_all.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **randomise_numberofpermutation** – Define the number of permutations. default=5000
- **skeletonised** – If True, randomize will be using only the white matter skeleton instead of the whole brain. default=True
- **metrics_dic** – Dictionnary containing the diffusion metrics to register in a common space. For each diffusion metric, the metric name is the key and the metric's folder is the value. default={'_noddi_odi': 'noddi', '_mf_fvf_tot': 'mf', '_diamond_kappa': 'diamond' }
- **regionWiseMean** – If true, csv containing atlas-based region wise mean will be generated.
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

regall(*folder_path=None, grp1=None, grp2=None, metrics_dic={'_diamond_kappa': 'diamond', '_mf_fvf_tot': 'mf', '_noddi_odi': 'noddi'}, slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Register all the subjects diffusion metrics specified in the argument metrics_dic into a common space using the transformation computed for the FA with the regall_FA function. This is performed based on TBSS of FSL. It is mandatory to have performed regall_FA prior to regall.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.

- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **metrics_dic** – Dictionary containing the diffusion metrics to register in a common space. For each diffusion metric, the metric name is the key and the metric's folder is the value. default={'_noddi_odi': 'noddi', '_mf_fvf_tot': 'mf', '_diamond_kappa': 'diamond'}
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slurm, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

regall_FA(*folder_path=None, grp1=None, grp2=None, starting_state=None, registration_type='-T', postreg_type='-S', prestats_treshold=0.2, slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Register all the subjects Fractional Anisotropy into a common space, skeletonised and non skeletonised. This is performed based on TBSS of FSL. It is mandatory to have performed DTI prior to regall_FA.

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.
- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **starting_state** – Manually set which step of TBSS to execute first. Could either be None, reg, post_reg, prestats, design or randomise. default=None
- **registration_type** – Define the argument used by the tbss command tbss_2_reg. Could either be '-T', '-t' or '-n'. If '-T' is used, a FMRIB58_FA standard-space image is used. If '-t' is used, a custom image is used. If '-n' is used, every FA image is align to every other one, identify the “most representative” one, and use this as the target image.
- **postreg_type** – Define the argument used by the tbss command tbss_3_postreg. Could either be '-S' or '-T'. If you wish to use the FMRIB58_FA mean FA image and its derived skeleton, instead of the mean of your subjects in the study, use the '-T' option. Otherwise, use the '-S' option.
- **prestats_treshold** – Thresholds the mean FA skeleton image at the chosen threshold during prestats. default=0.2
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 8 by a custom number of cpus of using slurm, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

tbss(*folder_path=None, grp1=None, grp2=None, starting_state=None, last_state=None, registration_type='-T', postreg_type='-S', prestats_threshold=0.2, randomise_numberofpermutation=5000, slurm=None, slurm_email=None, slurm_timeout=None, slurm_tasks=None, slurm_mem=None*)

Performs tract base spatial statistics (TBSS) between the data in *grp1* and *grp2*. The data type of each subject is specified by the *subj_type.json* file generated during the call to the *patient_list* function. The data type corresponds to the original directory of the subject (e.g. a subject that was originally in the folder *data_2* is of type 2). It is mandatory to have performed DTI prior to *tbss*. This function should not be used as it has been replaced by *regall_FA*, *regall* and *randomise_all* to allow for more flexibility.

example : `study.tbss(grp1=[1,2], grp2=[3,4])`

Parameters

- **folder_path** – the path to the root directory. default=*study_folder*
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.
- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **starting_state** – Manually set which step of TBSS to execute first. Could either be *None*, *reg*, *post_reg*, *prestats*, *design* or *randomise*. default=*None*
- **last_state** – Manually set which step of TBSS to execute last. Could either be *None*, *preproc*, *reg*, *post_reg*, *prestats*, *design* or *randomise*. default=*None*
- **registration_type** – Define the argument used by the *tbss* command *tbss_2_reg*. Could either be *'-T'*, *'-t'* or *'-n'*. If *'-T'* is used, a FMRIB58_FA standard-space image is used. If *'-t'* is used, a custom image is used. If *'-n'* is used, every FA image is align to every other one, identify the “most representative” one, and use this as the target image.
- **postreg_type** – Define the argument used by the *tbss* command *tbss_3_postreg*. Could either be *'-S'* or *'-T'*. If you wish to use the FMRIB58_FA mean FA image and its derived skeleton, instead of the mean of your subjects in the study, use the *'-T'* option. Otherwise, use the *'-S'* option.
- **prestats_threshold** – Thresholds the mean FA skeleton image at the chosen threshold during *prestats*. default=0.2
- **randomise_numberofpermutation** – Define the number of permutations. default=5000
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=*value_during_init*
- **slurm_email** – Email adress to send notification if a task fails. default=*None*
- **slurm_timeout** – Replace the default slurm timeout of 20h by a custom timeout.
- **slurm_tasks** – Replace the default number of slurm cpus of 8 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

white_mask(*folder_path=None, patient_list_m=None, corr_gibbs=True, forceUsePowerMap=False, debug=False, slurm=None, slurm_email=None, slurm_timeout=None, cpus=None, slurm_mem=None*)

Computes a white matter mask for each subject based on the T1 structural images or on the anisotropic power maps (obtained from the diffusion images) if the T1 images are not available. The outputs are available in the directories *<folder_path>/subjects/<subjects_ID>/masks/*. The T1 images can be gibbs ringing corrected.

example : `study.white_mask()`

Parameters

- **folder_path** – the path to the root directory. default=study_folder
- **patient_list_m** – Define a subset of subjects to process instead of all the available subjects. example : ['patientID1','patientID2','patientID3']. default=None
- **corr_gibbs** – If true, Gibbs ringing correction is performed on the T1 images. default=True
- **forceUsePowerMap** – Force the use of an AnisotropicPower map for the white matter mask generation. default=False
- **debug** – If true, additional intermediate output will be saved. default=False
- **slurm** – Whether to use the Slurm Workload Manager or not (for computer clusters). default=value_during_init
- **slurm_email** – Email adress to send notification if a task fails. default=None
- **slurm_timeout** – Replace the default slurm timeout of 3h by a custom timeout.
- **cpus** – Replace the default number of slurm cpus of 1 by a custom number of cpus of using slum, or for standard processing, its the number of core available for processing.
- **slurm_mem** – Replace the default amount of ram allocated to the slurm task (8096MO by cpu) by a custom amount of ram.

`elikopy.core.dicom_to_nifti(folder_path)`

Convert dicom data into compressed nifti. Converted dicoms are then moved to a sub-folder named original_data. The niftis are named patientID_ProtocolName_SequenceName.

Parameters **folder_path** – Path to root folder containing all the dicoms

12.3 elicopy.individual_subject_processing module

`elikopy.individual_subject_processing.diamond_solo(folder_path, p, core_count=4, reportOnly=False)`

Computes the DIAMOND metrics for a single subject. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/dMRI/microstructure/diamond/.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.
- **core_count** – Number of allocated cpu cores. default=4

`elikopy.individual_subject_processing.dti_solo(folder_path, p)`

Computes the DTI metrics for a single subject. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/dMRI/dti/.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.

`elikopy.individual_subject_processing.mf_solo(folder_path, p, dictionary_path, CSD_bvalue=None, core_count=1)`

Perform microstructure fingerprinting and store the data in the <folder_path>/subjects/<subjects_ID>/dMRI/microstructure/mf/.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.
- **dictionary_path** – Path to the dictionary of fingerprints (mandatory).
- **CSD_bvalue** – If the DIAMOND outputs are not available, the fascicles directions are estimated using a CSD with the images at the b-values specified in this argument. default=None
- **core_count** – Define the number of available core. default=1

`elikopy.individual_subject_processing.noddi_amico_solo(folder_path, p)`

Perform noddi amico on a single subject and store the data in the <folder_path>/subjects/<subjects_ID>/dMRI/microstructure/noddi_amico/.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.

`elikopy.individual_subject_processing.noddi_solo(folder_path, p, force_brain_mask=False, lambda_iso_diff=3e-09, lambda_par_diff=1.7e-09, use_amico=False, core_count=1)`

Computes the NODDI metrics for a single. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/dMRI/microstructure/noddi/.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.
- **force_brain_mask** – Force the use of a brain mask even if a whitematter mask exist. default=False
- **lambda_iso_diff** – Define the noddi lambda_iso_diff parameters. default=3.e-9
- **lambda_par_diff** – Define the noddi lambda_par_diff parameters. default=1.7e-9
- **use_amico** – If true, use the amico optimizer. default=FALSE
- **core_count** – Number of allocated cpu cores. default=1

`elikopy.individual_subject_processing.preproc_solo(folder_path, p, reslice=False, reslice_addSlice=False, denoising=False, gibbs=False, topup=False, topupConfig=None, forceSynb0DisCo=False, useGPUSynb0DisCo=False, eddy=False, biasfield=False, biasfield_bsplineFitting=[100, 3], biasfield_convergence=[1000, 0.001], starting_state=None, bet_median_radius=2, bet_numpass=1, bet_dilate=2, cuda=False, cuda_name='eddy_cuda10.1', s2v=[0, 5, 1, 'trilinear'], olrep=[False, 4, 250, 'sw'], qc_reg=True, core_count=1, niter=5, report=True, slspec_gc_path=None)`

Performs data preprocessing on a single subject. By default only the brain extraction is enabled. Optional preprocessing steps include : reslicing, denoising, gibbs ringing correction, susceptibility field estimation, EC-induced distortions and motion correction, bias field correction. The results are stored in the preprocessing subfolder of the study subject <folder_path>/subjects/<subjects_ID>/dMRI/preproc.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.
- **reslice** – If true, data will be resliced with a new voxel resolution of 2*2*2. default=False
- **reslice_addSlice** – If true, an additional empty slice will be added to each volume (might be useful for motion correction if one slice is dropped during the acquisition and the user still wants to perform easily the slice-to-volume motion correction). default=False
- **denoising** – If true, MPPCA-denoising is performed on the data. default=False
- **gibbs** – If true, Gibbs ringing correction is performed. We do not advise to use this correction unless the data suffers from a lot of Gibbs ringing artifacts. default=False
- **topup** – If true, Topup will estimate the susceptibility induced distortions. These distortions are corrected at the same time as EC-induced distortions if eddy=True. In the absence of images acquired with a reverse phase encoding direction, a T1 structural image is required. default=False
- **topupConfig** – If not None, Topup will use additional parameters based on the supplied config file located at <topupConfig>. default=None
- **forceSynb0DisCo** – If true, Topup will always estimate the susceptibility field using the T1 structural image. default=False
- **useGPUsynb0DisCo** – If true, Topup will estimate the susceptibility field with the T1 structural image using cuda. default=False
- **eddy** – If true, Eddy corrects the EC-induced (+ susceptibility, if estimated) distortions and the motion. If these corrections are performed the acquparam and index files are required (see documentation). To perform the slice-to-volume motion correction the slspec file is also needed. default=False
- **biasfield** – If true, low frequency intensity non-uniformity present in MRI image data known as a bias or gain field will be corrected. default=False
- **biasfield_bsplineFitting** – Define the initial mesh resolution in mm and the bspline order of the biasfield correction tool. default=[100,3]
- **biasfield_convergence** – Define the maximum number of iteration and the convergences threshold of the biasfield correction tool. default=[1000,0.001]
- **starting_state** – Manually set which step of the preprocessing to execute first. Could either be None, denoising, gibbs, topup, eddy, biasfield, report or post_report. default=None
- **bet_median_radius** – Radius (in voxels) of the applied median filter during brain extraction. default=2
- **bet_numpass** – Number of pass of the median filter during brain extraction. default=1
- **bet_dilate** – Number of iterations for binary dilation during brain extraction. default=2
- **cuda** – If true, eddy will run on cuda with the command name specified in cuda_name. default=False
- **cuda_name** – name of the eddy command to run when cuda==True. default="eddy_cuda10.1"
- **s2v** – list of parameters of Eddy for slice-to-volume motion correction (see Eddy FSL documentation): [mporder,s2v_niter,s2v_lambda,s2v_interp]. The slice-to-volume motion correction is performed if mporder>0, cuda is used and a slspec file is provided during the patient_list command. default=[0,5,1,'trilinear']

- **olrep** – list of parameters of Eddy for outlier replacement (see Eddy FSL documentation): [repol,ol_nstd,ol_nvox,ol_type]. The outlier replacement is performed if repol==True. default=[False, 4, 250, 'sw']
- **qc_reg** – If true, the motion registration step of the quality control will be performed. We do not advise to use this argument as it increases the computation time. default=False
- **niter** – Define the number of iterations for eddy volume-to-volume. default=5
- **slspec_gc_path** – Path to the folder containing volume specific slice-specification for eddy. If not None, eddy motion correction with gradient cycling will be performed.
- **report** – If False, no quality report will be generated. default=True
- **core_count** – Number of allocated cpu cores. default=1

`elikopy.individual_subject_processing.report_solo(folder_path, patient_path, slices=None, short=False)`

Legacy report function.

Parameters

- **folder_path** – path to the root directory.
- **patient_path** – Name of the subject.
- **slices** – Add additional slices cut to specific volumes
- **short** – Only output raw data, preprocessed data and FA data.

`elikopy.individual_subject_processing.white_mask_solo(folder_path, p, corr_gibbs=True, core_count=1, forceUsePowerMap=False, debug=False)`

Computes a white matter mask for a single subject based on the T1 structural image or on the anisotropic power map (obtained from the diffusion images) if the T1 image is not available. The outputs are available in the directories <folder_path>/subjects/<subjects_ID>/masks/. The T1 images can be gibbs ringing corrected.

Parameters

- **folder_path** – the path to the root directory.
- **p** – The name of the patient.
- **corr_gibbs** – If true, Gibbs ringing correction is performed on the T1 image. default=True
- **core_count** – Number of allocated cpu cores. default=1
- **forceUsePowerMap** – Force the use of an AnisotropicPower map for the white matter mask generation. default=False
- **debug** – If true, additional intermediate output will be saved. default=False

12.4 elicopy.utils module

`elikopy.utils.anonymise_nifti(rootdir, anonymize_json, rename)`

Anonymise all nifti present in rootdir by removing the PatientName and PatientBirthDate (only month and day) in the json and renaming the nifti file name to the PatientID.

Parameters

- **rootdir** – Folder containing all the nifti to anonymise.

- **anonymize_json** – If true, edit the json to remove the PatientName and replace the Patient-BirthDate by the year of birth.
- **rename** – If true, rename the nifti to the PatientID.

`elikopy.utils.export_files(folder_path, step, patient_list_m=None)`

Creates an export folder in the root folder containing the results of 'step' for each patient in a single folder

example : `export_files('user/my_rootfolder', 'dMRI/microstructure/dti')`

Parameters

- **folder_path** – root folder
- **step** – step to export
- **patient_list_m** – Define a subset a patient to process instead of all the available subjects.

`elikopy.utils.getJobsState(folder_path, job_list, step_name)`

Periodically checks the status of all jobs in the `job_list`. When a job status change to complete or a failing state. Write the status in the log and remove the job from the `job_list`. This function end when all jobs are completed or failed.

Parameters

- **folder_path** – The path to the root dir of the study (used to write the logs.txt file)
- **job_list** – The list of job to check for state update
- **step_name** – The string value of the prefix to put in the log file

`elikopy.utils.get_job_state(job_id)`

Retrieve the state of a job through the `sacct` bash command offered by the `lurm` Workload Manager. :param `job_id`: The id of the job to retrieve the state of. :return state: The string value representing the state of the job.

`elikopy.utils.get_patient_list_by_types(folder_path, type=None)`

Print the list of patient corresponding to a specific type of patient.

Parameters

- **folder_path** – Path to the root folder of the study.
- **type** – The selected type

`elikopy.utils.inference(T1_path, b0_d_path, model, device)`

`synb0DISCO` adapted from <https://github.com/MASILab/Synb0-DISCO>

Parameters

- **T1_path** – Path to the normalized projected T1.
- **b0_d_path** – Path to the b0 atlases.
- **model** – DL Model
- **device** – Define if cuda or cpu is used.

`elikopy.utils.makedir(dir_path, log_path, log_prefix)`

Create a directory in the location specified by the `dir_path` and write the log in the `log_path`.

Parameters

- **dir_path** – The path to the directory to create.
- **log_path** – The path to the log file to write verbose data.
- **log_prefix** – The prefix to use in the log file.

`elikopy.utils.merge_all_reports(folder_path)`

Merge all subjects quality control reports into a single report.

Parameters `folder_path` – Path to the root folder of the study.

`elikopy.utils.merge_all_specific_reports(folder_path, merge_wm_report=False, merge_legacy_report=False)`

Merge all selected specific subject's report into a single big report.

Parameters

- **folder_path** – Path to the root folder of the study.
- **merge_wm_report** – Select wm report.
- **merge_legacy_report** – Select legacy report.

`elikopy.utils.randomise_all(folder_path, randomise_numberofpermutation=5000, skeletonised=True, metrics_dic={'FA': 'dti', '_diamond_kappa': 'diamond', '_mf_fvf_tot': 'mf', '_noddi_odi': 'noddi'}, core_count=1, regionWiseMean=True)`

Performs tract base spatial statistics (TBSS) between the data in grp1 and grp2 (groups are specified during the call to `regall_FA`) for each diffusion metric specified in the argument `metrics_dic`. The mean value of the diffusion metrics across atlases regions can also be reported in CSV files using the `regionWiseMean` flag. The used atlases are : the Harvard-Oxford cortical and subcortical structural atlases, the JHU DTI-based white-matter atlases and the MNI structural atlas It is mandatory to have performed `regall_FA` prior to `randomise_all`.

Parameters

- **folder_path** – path to the root directory.
- **randomise_numberofpermutation** – Define the number of permutations. default=5000
- **skeletonised** – If True, randomize will be using only the white matter skeleton instead of the whole brain. default=True
- **metrics_dic** – Dictionnary containing the diffusion metrics to register in a common space. For each diffusion metric, the metric name is the key and the metric's folder is the value. default={'_noddi_odi': 'noddi', '_mf_fvf_tot': 'mf', '_diamond_kappa': 'diamond'}
- **regionWiseMean** – If true, csv containing atlas-based region wise mean will be generated.
- **core_count** – Number of allocated cpu core. default=1

`elikopy.utils.regall(folder_path, grp1, grp2, core_count=1, metrics_dic={'_diamond_kappa': 'diamond', '_mf_fvf_tot': 'mf', '_noddi_odi': 'noddi'})`

Register all the subjects diffusion metrics specified in the argument `metrics_dic` into a common space using the transformation computed for the FA with the `regall_FA` function. This is performed based on TBSS of FSL. It is mandatory to have performed `regall_FA` prior to `regall`.

Parameters

- **folder_path** – path to the root directory.
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.
- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **metrics_dic** – Dictionnary containing the diffusion metrics to register in a common space. For each diffusion metric, the metric name is the key and the metric's folder is the value. default={'_noddi_odi': 'noddi', '_mf_fvf_tot': 'mf', '_diamond_kappa': 'diamond'}
- **core_count** – Define the number of available core. default=1

`elikopy.utils.regall_FA(folder_path, grp1, grp2, starting_state=None, registration_type='-T', postreg_type='-S', prestats_treshold=0.2, core_count=1)`

Register all the subjects Fractional Anisotropy into a common space, skeletonised and non skeletonised. This is performed based on TBSS of FSL. It is mandatory to have performed DTI prior to `regall_FA`.

Parameters

- **folder_path** – path to the root directory.
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.
- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **starting_state** – Manually set which step of TBSS to execute first. Could either be `None`, `reg`, `post_reg`, `prestats`, `design` or `randomise`. default=`None`
- **registration_type** – Define the argument used by the `tbss` command `tbss_2_reg`. Could either be `'-T'`, `'-t'` or `'-n'`. If `'-T'` is used, a FMRIB58_FA standard-space image is used. If `'-t'` is used, a custom image is used. If `'-n'` is used, every FA image is align to every other one, identify the “most representative” one, and use this as the target image.
- **postreg_type** – Define the argument used by the `tbss` command `tbss_3_postreg`. Could either be `'-S'` or `'-T'`. If you wish to use the FMRIB58_FA mean FA image and its derived skeleton, instead of the mean of your subjects in the study, use the `'-T'` option. Otherwise, use the `'-S'` option.
- **prestats_treshold** – Thresholds the mean FA skeleton image at the chosen threshold during `prestats`. default=0.2
- **core_count** – Define the number of available core. default=1

`elikopy.utils.submit_job(job_info)`

Submit a job to the Slurm Workload Manager using a crafted `sbatch`.

Parameters `job_info` – The parameters to use in the `sbatch`.

Return `job_id` The id of the submitted job.

`elikopy.utils.synb0DisCo(folder_path, topuppath, patient_path, starting_step=None, topup=True, gpu=True)`
`synb0DISCO` adapted from <https://github.com/MASILab/Synb0-DISCO>

Parameters

- **folder_path** – path to the root directory.
- **topuppath** – Path to the subject's topup folder.
- **patient_path** – Name of the subject.
- **starting_step** – Define the starting step, usefull if previous step had already been run.
- **topup** – If true, topup will be performed after `synb0DisCo`.
- **gpu** – If true, torch will use the gpu.

Return type object

`elikopy.utils.tbss_utils(folder_path, grp1, grp2, starting_state=None, last_state=None, registration_type='-T', postreg_type='-S', prestats_treshold=0.2, randomise_numberofpermutation=5000)`

[Legacy] Performs tract base spatial statistics (TBSS) between the data in `grp1` and `grp2`. The data type of each subject is specified by the `subj_type.json` file generated during the call to the `patient_list` function. The data type corresponds to the original directory of the subject (e.g. a subject that was originally in the folder `data_2` is of type 2). It is mandatory to have performed DTI prior to `tbss`.

Parameters

- **folder_path** – path to the root directory.
- **grp1** – List of number corresponding to the type of the subjects to put in the first group.
- **grp2** – List of number corresponding to the type of the subjects to put in the second group.
- **starting_state** – Manually set which step of TBSS to execute first. Could either be None, reg, post_reg, prestats, design or randomise. default=None
- **last_state** – Manually set which step of TBSS to execute last. Could either be None, preproc, reg, post_reg, prestats, design or randomise. default=None
- **registration_type** – Define the argument used by the tbss command tbss_2_reg. Could either be ‘-T’, ‘-t’ or ‘-n’. If ‘-T’ is used, a FMRIB58_FA standard-space image is used. If ‘-t’ is used, a custom image is used. If ‘-n’ is used, every FA image is align to every other one, identify the “most representative” one, and use this as the target image.
- **postreg_type** – Define the argument used by the tbss command tbss_3_postreg. Could either be ‘-S’ or ‘-T’. If you wish to use the FMRIB58_FA mean FA image and its derived skeleton, instead of the mean of your subjects in the study, use the ‘-T’ option. Otherwise, use the ‘-S’ option.
- **prestats_threshold** – Thresholds the mean FA skeleton image at the chosen threshold during prestats. default=0.2
- **randomise_numberofpermutation** – Define the number of permutations. default=5000

12.5 Module contents

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`tbss_utils()` (*in module elikopy.utils*), [52](#)

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