# mreg\_recon\_tool

Manual for the MREG data reconstruction Tool. The interested reader can find additional background information on the topic of reconstructing MREG data in the following by Thimo Hugger dissertation (only in german): http://www.freidok.uni-freiburg.de/volltexte/8081/

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In the manual the so called MREG-sequence is mentioned several times, by this either the "MREG\_stable", the "MREG\_c2p" or the "MREG\_c2p\_VD11" sequences are meant which were implemented by Benjamin Zahneisen and Jakob Assländer.

# 1. Top Elements

### 1.1. Push button: Set Working Dir

This let's you select a directory, that is used as default for saving any data. By default an additional folder is created in this folder which then contains the reconstructed data. The currently active working directory is displayed in the edit box next to the push button.

# 1.2. Push button: Export Settings

Let's you save all the settings in the current mreg\_recon\_tool to a .mat-file. You can then open a new tool and provide the exported data as an argument to mreg\_recon\_tool. It will then automatically try to reach the state of the tool that was used to export the settings at the time of the export.

### 2. Raw Data

#### 2.1. Push button: Load RawData

Here the file that contains the raw data can be selected. The file needs to be a .dat-file that was measured with the MREG-sequence. The callback of the push button calls the function loadData.m. The file is not actually loaded, instead only the header information is read. Later on each reconstruction of a timeframe then only loads the necessary part from the raw data,

by calling loadData(filename, timeframe). The push button labeled X next to it clears the currently loaded raw data.

# 3. Reference

#### 3.1. Push button: Load Reference

Let's you select the reference data file. It can either be a .dat-file or a Matlab .mat-file, where the .mat-file is the reference data saved with the save reference button 3.4. If the selected file is a .dat-file, the callback executes loadReference(filename, sensmode), where sensmode is given by the popup menu *Sens. mode* 3.2.

### 3.2. Popup menu: Sensitivity mode

The mode to compute the sensitivity maps from the reference data can be set here. The mode is used when the *Load Reference* button is pressed. In case the reference data has already been loaded, the sensitivity maps will be recalculated with the selected mode, using the function coilSensitivities. The following modes are available:

- 1. *sos*: The coil sensitivity of a coil is equal to the single coil image divided by a sum of squares image of all coil images.
- 2. *adapt*: The coil sensitivities are computed using the function adapt3D. Each coil image is divided by an optimized complex combination of all coil images.
- 3. *lowres*: The coil sensitivity is given by a strongly low-pass filtered version of the coil image.

### 3.3. Popup menu: Select view

Determines what will be displayed when the view button is pressed. The following options exist:

- 1. Coil Profiles: Reference images of each coil.
- 2. *Coil Sensitivities*: The computed coil sensitivities.
- 3. Off-resonance map: A map of the estimated off-resonances due to  $B_0$  field inhomogeneities in radians per second.

### 3.4. Push button: Save

Allows to save the data as a .mat-file. Since loading the data from a .dat-file can be time consuming, saving the data to a .mat-file can speed up the loading process a lot for future sessions.

#### 3.5. Push button: View

The data selected in 3.3 is displayed using the viewer3d\_tool.

# 4. Trajectory

### 4.1. Push button: Load Trajectory

A trajectory may be loaded from a file, where the file must either be a series of .dat-files, a .mat-file or a .grad-file .dat-files correspond to a set of trajectory calibration measurements, acquired with the MREG-sequence.

Three .dat-files must be selected from the uigetfile dialogue (use the CTRL-key to select more than one file), corresponding to separate measurements of the three components (labeled "phase,," "read,, and "slice,, in the MREG-sequence) of the trajectory.

A .grad-file can also be used in case a measurement of the trajectory is not available.

Furthermore the trajectory can be contained in a .mat-file, that was created using the save button 4.3.

The callback will then execute the function loadTrajectory\_MatDatGrad which handles all three data types.

If the .grad file is stored either in the same folder as the raw data or in (mat-lab\_new/fmri/)mreg/grad\_files, the trajectory is loaded automatically. The latter folder is meant to store all common trajectories.

### 4.2. Push button: Select range

Allows you to select only part of the trajectory. Most of the time one will want to discard points in the beginning and at the end of the data that corresponds to parts of the trajectory that are either unimportant or might affect the reconstruction quality negatively (e.g. parts where the trajectory does not change). Usually the optimal range is already read from the header of the .grad or the .dat file, so no further selection is necessary.

If the button is pressed, a window will pop up, showing the length of the k-spaces point versus the index, as guidance for the selection. One can then click on two positions in the graph and the range between these two points will be used, while the rest will be discarded. The process

will repeat in case there is more than one interleaf.

#### 4.3. Push button: Save

Allows to save the trajectory to a .mat-file. The .mat-file will also contain the range information, so that saving the trajectory allows to skip the range selection step in future sessions.

#### 4.4. Push button: View

If this button is pressed the trajectory will be displayed in a plot.

### 5. Reconstruction Parameters

#### 5.1. Editbox: CG-Tolerance

The value entered here will determine when the conjugate (nonlinear) gradient iterations will be stopped. A value around 1e-5 is usually reasonable.

#### 5.2. Editbox: Max. Iterations

The value entered here will determine the maximum number of iterations the (nonlinear) conjugate gradient algorithm will perform before it is stopped. If only the  $L_2$ -norm is used as a penalty in 7.3 then the standard conjugate gradient algorithm will be employed and a reasonable value for the maximum number of iterations is around 40. In case the  $L_1$ -norm is used the nonlinear conjugate gradient algorithm is used. Due to the slower convergence a value of around 200 should be used here.

#### 5.3. Editbox: Recon Size

The array size of the reconstruction can be adjusted here. This option as well as the next (Voxel Size) will only be enabled once the reference and the trajectory are loaded. Using a smaller reconstruction size can greatly benefit the reconstruction time since a smaller array size speeds up the internal operations like e.g. the FFT. Please note that this will not affect the resolution, that means reducing the array size leads to a reduction of the field of view. One therefore needs to make sure that the object still fits into the reduced field of view, otherwise aliasing will occur.

### 5.4. Editbox: Voxel Size

If really necessary the effective voxel size in the reconstruction can be changed here. Internally this will lead to the application of scaling factors to the k-space trajectory. This can be used e.g. to reduce the array size of the reconstruction beyond the point where the object would be larger than the field of view, by enlarging the voxel size.

### 5.5. Push button: Adjust

In case the trajectory and the reference have different size and resolution and in case the Recon Size or the Voxel Size are changed, one has to adjust the reference data to the new settings. Internally the coil sensitivities, the coil profiles and the off-resonance map are resized. The callback mainly executes the function imresize3D to do that. If correctly adjusted, the color of the button will change to green.

# 6. Reconstruction Options

### 6.1. Popup menu: Reconstruction Type

This option allows to set the reconstruction method in case of an interleaved acquisiton<sup>1</sup> If the acquisition was interleaved then *sliding window* or *KWIC* need to be used to get consistent results within timeframes, otherwise *standard* is usually the correct choice.

- 1. *standard*: Reconstructs each timeframe separately using the corresponding interleaf of that timeframe.
- 2. sliding window: If  $N_i$  interleaves are used, and timeframe n is to be reconstructed, the neighboring timeframes  $(n-{\sf floor}(\frac{N_i}{2}+1)+1)\dots(n+Ni-{\sf floor}(\frac{N_i}{2}+1))$  are combined and treated as the current timeframe. If the neighboring timeframe exceed the range of the available timeframes then the reconstruction of the current timeframe is set to zero.
- 3. *KWIC*: Similar to the previous option. Differences are that less data is used from interleaves that are further away from the current timeframe and the timeframes used are given by  $n \dots n + N_i 1$ . This option is most sensible when the so called bit-reversed ordering is used for the interleaves.

<sup>&</sup>lt;sup>1</sup>If  $k = 1 ... N_i$  different trajectories where used for data acquisition this typically means that in timeframe with index n, k is equal to  $rem(n, N_i)$ .

### 6.2. Checkbox: Off-resonance Correction

If this option is checked an off-resonance effects will be incorporated in the image reconstruction by using the off-resonance map given by *Load Reference* 3.1. If this option is checked the reconstruction time per timeframe is prolonged approximately by a factor 10.

### 6.3. Checkbox: z0

When z0 is chosen, one image in the middle of the time series is reconstructed. This image is then used as the starting point for the conjugate gradient. The number of steps necessary for convergence should be smaller than when using zero as starting point (default). But this option should be handled with care, since too few steps could give you a image, that corresponds more to the initial guess than to the image at that particular time frame. This could result in a meaningless time series.

### 6.4. Checkbox: DORK

"Dynamic off-resonance in k-space" is a method by Josef Pfeuffer et al. (MRM, 2002), that corrects for  $B_0$  changes due to scanner drifts, respiration etc. They way it works: If the frequency is off-center, the phase of the FID develops over time. If the the change of phase changes over the time series, the global off-resonance frequency can be calculated for the particular time frame and multiplied to the raw data. But it assumes, that the first time frame is on resonance. So I highly recommend to adjust the frequency before every MREG scan (Options -> Adjustments in the exam task of the scanner).

The computational costs are mainly a couple of seconds when turning on the mode. This is when the frequency drift is calculated. For the reconstruction itself the costs are negligible. Therefore I highly recommend to use the option.

It seems to work very well for the drift of the  $B_0$  field, making sure, that last image is not shifted compared to the first and is of reasonable quality. The respiration induced artifacts seem to change not that much, even though they are clearly visible in the frequency (have a look in recon\_details.dork\_frequency either stored by the grid engine or as a variable in mreg\_recon\_tool\_recon; just set a breakpoint and plot the vector). The reason, why the respiration artifacts are not reduced by that much are higher spatial order components of the frequency (see Magnetic Resonance Encephalography Reconstruction with Magnetic Field Monitoring, F. Testud et al., ISMRM 2012).

### 6.5. Editbox: Global off-resonance

Global means in this case "4D-global". So, if you forgot to adjust the frequency before the MREG scan, the frequency can be wrong right from the beginning of the time series, which cannot be corrected by DORK. This will result in crappy images with lots of shading all over the brain. If that is the case, you can try to find a frequency that improves the image. Just try some values, usually between -100/s and 100/s and see which one fits best. There is no, automatic or systematic way for that, sorry.

### 6.6. Checkbox: Use GridEngine

If the Sun GridEngine (http://wikis.sun.com/display/GridEngine/Home) is available to you and you want to reconstruct a lot of timeframes, this option can be checked to split up the reconstruction task into jobs and distribute them among the available slots on the GridEngine. Depending on the available resources this can tremendously speed things up. The option will be grayed out if the GridEngine is not found on your system.

### 6.7. Editbox: Frames per job

If the GridEngine option is checked, one can specify here the number of timeframes that will comprise one GridEngine job.

## 6.8. Editbox: Timeframe Range

Here you can specify the timeframes that you would like to reconstruct. Any vector of timeframes is acceptable as long as each timeframe is within the boundaries given by data. Standard Matlab syntax can be used in this editbox, like for example [101:2:200]. Please note that unless the data is saved to files during the reconstruction (see 6.10) the resulting 4D-array will contain no information about the timeframes that were used.

### 6.9. Push button: Start Recon

Performs the Reconstruction either locally or submits jobs to the gridengine depending on the option 6.6.

# 6.10. Checkbox: Save during recon

If this option is checked the data saved to the current working directory 1.1 in a folder with the name given by the option *Where* 6.12 if no name is specified or *Where* is empty then a

default folder will be created containing the current timestamp to avoid overwriting of previous reconstructions. This option will become unavailable if *Use GridEngine* 6.6 is checked, since then each job saves the data during reconstruction by default.

### 6.11. Popup menu: Save as ...

Sets the file format for the saved reconstructions. Each frame will be saved in a separate file. The filename will contain the timefame of the reconstruction in the file.

- 1. *mat*: Standard Matlab mat-file. Contains no additional information about the reconstruction at this point.
- 2. *nifit*: NIFTI file format. The header of the resulting .nii-file contains some information about the reconstruction. Please note that SPM is needed in order to save to this file format.

#### 6.12. Push button: Where

Optional possibility to name the folder inside the working directory 1.1 where the data will be saved. This is useful in case you want to open several mreg\_recon\_tools to split up a local reconstruction task in several smaller tasks and make sure they all are saved in the same folder. If this is empty a default folder will be created containing a timestamp.

#### 6.13. Push button: Save File

If *Save during recon*. was not checked and a reconstruction was performed it will be viewable only inside the mreg\_recon\_tool. This option allows to save the data to a file in the file format given by 6.11.

### 6.14. Push button: Save WS

Sometimes one might want to analyze the reconstruction inside the Matlab workspace. Use this option to export the data to the current Matlab workspace. A window will then pop up, where you can name the variable.

#### 6.15. Push button: View

If reconstructed data is available opens the viewer3d\_tool to view it.

### 7. Penalties

### 7.1. Push buttons: Add & Delete

Adds and deletes penalties in the cost function. The default cost function always consists of the fidelity term. If A is the encoding operator,  $\rho$  is the unknown image and b is the measured data, then the cost function has the form

$$f(\boldsymbol{\rho}) = \|A\boldsymbol{\rho} - \boldsymbol{b}\|_2^2.$$

Adding a penalty (pressing the Push button Add), really just adds the specified penalty mathematically to the fidelity term. With e.g. a regularization parameter  $\lambda$ , the L<sub>1</sub>-norm and an operator W the resulting cost function would be equal to:

$$f(\boldsymbol{\rho}) = \|A\boldsymbol{\rho} - \boldsymbol{b}\|_2^2 + \lambda \|W\boldsymbol{\rho}\|_1.$$

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### 7.2. Editbox: lambda

The regularization parameter for this penalty.

# 7.3. Popup menu: Select norm

The norm for the penalty can be selected here. It can either be the  $L_1$ - or the  $L_2$ -norm.

# 7.4. Popup menu: Select operator

An operator can be selected that is applied to  $\rho$  before measuring the norm of it.