# QTI+: MATLAB Implementation User Guide

# Contents

1	Introduction			
	1.1	Software overview	2	
2	Insta 2.1 2.2	allation CVX		
3	J			
	3.1	Input: Data & B-tensor format	5	
	3.2	qtiplus_fit		
		3.2.1 choosing the pipeline	6	
		3.2.2 passing a mask		
		3.2.3 choosing the SDP solver	7	
		3.2.4 excluding measurements from the fit	7	
		3.2.5 Using the parallel computing toolbox	8	
		3.2.6 selecting the number of voxels to be fitted simultaneously	8	
	3.3	Output: model parameters and scalar invariants	9	
	3.4	Computational times	9	
4	Exar	mple	10	

## 1 Introduction

This document contains the User Guide for the QTI+ software as implemented in Matlab (https://github.com/DenebBoito/qtiplus). The document is structured in three sections. The first contains instructions on the necessary modules that need to be installed and how to install them. Then comes the main section describing how to process diffusion MRI data acquired with time-varying magnetic field gradients using this software. The final section contains an example with code snippets showing how to process an example dataset using different options.

This guide assumes that the reader is or will be familiar with the articles describing the QTI framework [1], linked here, and the methods belonging to the "+" version, linked here. The focus of this guide lies only on providing an overview of the software functionalities and how to use them. For the terminology, methods, and science, we refer the reader to the articles.

#### 1.1 Software overview

The QTI+ framework and the nomenclature are displayed in Figure 1. The different blocks correspond to the different fitting routines as presented in [2], please refer to that paper for information regarding the various steps.

For the SDPdc, (m) check, and SDPdcm steps, one of the solvers interfaced by CVX is used. CVX is introduced in section 2.1, while information about the solvers are found in section 3.2.3. For solving the nonlinear problem, the Matlab routine *lsqcurvefit* with the *Levenberg-Marquardt* algorithm is used. Note that, since the solution from the SDPdc step is used as starting point for the nonlinear fit, if the residuals after the SDPdc step are smaller than those after the NLLSdc step, the solution produced by SDPdc is retained.

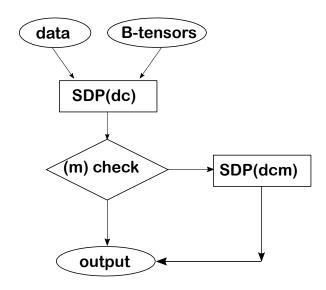
Note that it is not necessary to perform all the steps in the framework; satisfactory results can be obtained just by performing SDPdc. In the software, the user can decide which of the steps to perform by selecting predefined pipelines (By default, only SDPdc is performed). The main output consists of the estimated QTI model parameters. Additionally, the user can obtained the scalar maps derived from these estimates.

The software is really intended to be used through the main function called <a href="mailto:qtiplus\_fit">qtiplus\_fit</a>. This is the function that takes the data and B-tensors as inputs, and returns the estimated model parameters and the scalar invariants derived from these estimates. Details about the expected format for the data and b-tensors are given in section 3.1. Used in the following fashion:

```
[model, invariants] = qtiplus_fit(data,btensors);
```

the software uses all the default settings, which is possibly what most users would like to do. Details about the different settings can be found in sections 3.2.1, 3.2.2, 3.2.3, 3.2.4, and 3.2.5

#### **QTI+ Flowchart**



#### Nomenclature:

SDPdc: Semidefinite Programming formulation with (d) and (c) constraints

NLLSdc: Non-linear least square formulation with (d) and (c) constraints

(m) check: Check on condition (m) via Semidefinite Programming

SDPdcm: Semidefinite Programming formulation with (d), (c), and (m) constraints

Figure 1: QTI+ framework layout and nomenclature.

## 2 Installation

#### 2.1 CVX

In order to use the routines in this library, CVX (http://cvxr.com/cvx/) needs to be installed [3]. CVX offers an environment for defining and solving convex problems. It is free to download and install, and comes with a few free solvers that can be readily used upon installation. If CVX is not already installed, please visit the download page (http://cvxr.com/cvx/download/) and download the option that suits you. Please consult the installation page (http://cvxr.com/cvx/doc/install.html) for instruction on how to install CVX on your machine.

Note that in our implementation we favoured Mosek (MOSEK ApS, Denmark) as the solver to be used for solving the problem using Semidefinite Programming (SDP). Even though Mosek can be obtained together with CVX (depending on which bundle is downloaded), it requires a licence in order to be used. Such licence is free for academic purposes, and instructions on how to obtain the licence and where to place it in your machine can be found here

http://cvxr.com/cvx/doc/mosek.html. If Mosek is not available, the QTI+ routines will use another solver that ships with CVX, SDPT3 [4]. There should be no difference in terms of results when choosing one or the other solver. In our experience, the only relevant difference is found in the computational times, with Mosek being the faster option. In section 3.4 you may find a comparison between Mosek and SDPT3 in terms of computational times.

## 2.2 QTI+

All the codes for QTI+ are written in Matlab, meaning that it is sufficient to either clone or download the files from the Github page, and add them to your Matlab path to start using it. The function <a href="mailto:qtiplus\_setup">qtiplus\_setup</a> may assist in adding all the folders to the path, and should be run before anything else to avoid error messages. This function is also called when running the script containing the example on how to use the software example\_qtiplus\_fit.

# 3 Processing the data

In this section we detail inputs, outputs, and optional parameters that can be set in the software.

# 3.1 Input: Data & B-tensor format

The data is expected to be a 4D Matlab matrix with shape [nx, ny, nz, nd]. nx, ny, and nz are the number n of voxels for each dimension x, y, and z. The nd different diffusion volumes are stacked along the 4th dimension. The example dataset provided with the package in the example dataset folder has dimensions [64, 64, 6, 217]. This means, there are 6 xy-plane slices of size  $64 \times 64$ , and there are 217 diffusion measurements.

The b-tensors are expected to be shaped as a 2D Matlab matrix with shape [nd, 6], where the number of rows represents the number of collected diffusion volumes, and each row contains one b-tensor in Voigt format. For the example mentioned above, the matrix containing all the B-tensors has size [217, 6]. The following convention for converting a B-tensor from a  $3 \times 3$  matrix representation to its  $1 \times 6$  Voigt representation is adopted:

$$\begin{pmatrix} B_{xx} & B_{xy} & B_{xz} \\ B_{xy} & B_{yy} & B_{yz} \\ B_{xz} & B_{yz} & B_{zz} \end{pmatrix} \longleftrightarrow \begin{bmatrix} B_{xx} & B_{xy} & B_{xz} & \sqrt{2}B_{xy} & \sqrt{2}B_{xz} & \sqrt{2}B_{yz} \end{bmatrix}$$

The functions named convert\_\* in the helper functions folder may assist the user in switching from one representation to the other.

Please note that the B-tensors are expected to have SI units, i.e.  $s/m^2$ !

# 3.2 qtiplus\_fit

This is the main function of the package which is intended as interface between the user and the software. It controls the input, output, and behaviour of the fit via different key-value pairs. Concise information about input, output, and optional settings can be obtained by typing "help qtiplus\_fit" in the Matlab command window or by looking at the comments in the script defining the function. In the next sections the available options for controlling the software behaviour via key-value pairs are detailed. For each controllable option, the keyword, the available options, and the default value are highlighted in a box. The accompanying text provides more insights on choosing between the different options.

## 3.2.1 choosing the pipeline

```
KEY: 'pipeline'

VALUE OPTIONS:

0: SDPdc

1: SDPdc → NLLSdc

2: SDPdc → NLLSdc → (m)-check → SDPdcm

3: SDPdc → NLLSdc → SDPdcm

4: SDPdc → (m)check → SDPdcm

5: SDPdc → SDPdcm

DEFAULT: 0
```

The keywoard 'pipeline' allows the user to select which of the steps in the QTI+ framework to perform. As mentioned in the introduction, it is not necessary to perform all of the steps in the framework as satisfactory results can be obtained employing SDPdc only. This saves computational time and is therefore the default options. For a comparison between computational times using different options, the user is referred to [2] and to section 3.4.

The choice between different pipelines is made by passing an integer after the keyword. For example, to analyse the data using all the steps in the QTI+ framework, which corresponds to option 2 as shown in the box, the qtiplus\_fit function can thus be called in the following way:

```
[model, invariants] = qtiplus_fit(data,btensors, 'pipeline', 2);
```

#### 3.2.2 passing a mask

KEY: 'mask'

VALUE OPTIONS: 3D binary image volume

**DEFAULT:** if none is input, one is computed using the median\_otsu method implemented similarly to that in Dipy [5].

The purpose of passing a mask is to avoid computations in parts of the image that do not belong to the object, or to perfom computations only on selected regions of interest. For example, in neuroimaging it is quite common to do computations only on voxels containing brain tissue. The user has here the choice to either pass a mask to the qtiplus\_fit function, or let the software compute one using the median\_otsu method implemented similarly to that in Dipy [5]. The following call shows how to pass a mask to the fitting function. The variable "mymask" contains the 3D matrix with binary (or logical) values, and has the same [nx, ny, nz] size as the input data.

```
[model, invariants] = qtiplus_fit(data,btensors, 'mask', mymask);
```

Note: if it is desired to compute the model parameters on all parts of the volume, it is recommended that the corresponding mask is input to the <a href="mailto:qtiplus\_fit">qtiplus\_fit</a> function. This would correspond to the following call:

```
mymask = ones(nx, ny, nz);
[model, invariants] = qtiplus_fit(data,btensors, 'mask', mymask);
```

Note: The implemented method for automatically computing a mask works fine on the datasets we have tried it on, but it is not guaranteed that it will work fine on all possible datasets. The user should therefore always check that the automatically computed mask is ok. This can be verified by checking the output of the simple\_mask function located in the helper functions. For the data that will be analysed, the following call will return the mask that would be automatically computed if none is passed to the <a href="mailto:qtiplus\_fit">qtiplus\_fit</a>:

```
mymask = simple_mask(data);
```

#### 3.2.3 choosing the SDP solver

```
KEY: 'solver'
VALUE OPTIONS:
    'mosek'
    'sdpt3'

DEFAULT: 'mosek' if available, otherwise revert to 'sdpt3'
```

This key-value pair determines which of the solvers to use for solving the SDP problem. If no option is input, the function <code>qtiplus\_fit</code> checks whether Mosek is available, and if so, it uses it. If Mosek is not available, the software reverts to SDPT3, which is installed together with CVX. Mosek is generally recommended compared to other solvers for convex programming, and it is the solver that was used to produce the results in [2]. In terms of results, there should not be any difference between Mosek and SDPT3. In terms of speed, Mosek is in our experience the faster option. In section 3.4 there is a comparison between the computational times achievable with both Mosek and SDPT3 on the example dataset provided with this software package.

#### 3.2.4 excluding measurements from the fit

KEY: 'ind'

VALUE OPTIONS: binary array of size [nd,1] indicating which measurements to consider for

the fitting.

**DEFAULT:** all measurement points are used

By using this option the user can select to exclude some of the collected data from being part of the fitting. For example, in diffusion MRI, when applying a model to some data, it is quite common to exclude the measurements acquired without diffusion encoding gradients applied. Considering the example dataset provided with the software package, the first 13 measurements were acquired without diffusion weighting. Therefore we can choose to exclude them from the fitting. This is done in the following way: first, we define an array with zeros on the indices of the measurements that we want to exclude, and ones on the indices of the measurements that we want to keep. In our example, this corresponds to an array with zeros

in the first 13 positions, and ones everywhere else:

```
indx = ones(217,1);
indx(1:13) = 0;
```

Then, we pass this array to the <a href="mailto:qtiplus\_fit">qtiplus\_fit</a> function with the keyword 'ind'.

```
[model, invariants] = qtiplus_fit(data, btensors, 'ind', indx);
```

#### 3.2.5 Using the parallel computing toolbox

KEY: 'parallel'

VALUE OPTIONS:

0: single worker computation

1: computation distributed over multiple workers

**DEFAULT: 1** 

The QTI+ fitting routines can take advantage of computations being independent between voxels, thus meaning that they can be performed in parallel. In Matlab, distributing the operations over multiple workers requires the *Parallel Computing Toolbox*. If the option for performing parallel computations is selected, the software first check whether this toolbox is available, and if so exploits as many workers as per Matlab's settings. If the user prefer instead to not perform computations in parallel, this needs to be enforced in the following way:

```
[model, invariants] = qtiplus_fit(data, btensors, 'parallel', 0);
```

## 3.2.6 selecting the number of voxels to be fitted simultaneously

KEY: 'nvox'

VALUE OPTIONS: integer indicating how many voxels to be fitted simultaneously.

**DEFAULT:** 50

As there is no dependence between the fits performed in each voxel, several can be fitted independently simultaneously. This parameter thus represents the number of voxels that get fitted at the same time. This applies to SDPdc, (m)-check, and SDPdcm. The step involving the nonlinear fit, NLLSdc, can only be applied on one voxel at a time.

Note the difference between this and the performing parallel computations: performing the computations on multiple workers means that each worker gets to solve one problem. With the *nvox* parameter we decide instead over how many voxels one problem is defined. Thus, for example, if 4 workers are available, and we select nvox = 50, 200 voxels are being fitted simultaneously.

Performing the fit with nvox > 1 provides a great speed up in computations, since the major bottleneck for SDPdc, (m) check, and SDPdcm is the repeated preconditioning done by the solvers before solving the given problem. By constructing a problem over several voxels, we

avoid having to do the preconditioning for each of them. Note however that there is a trade-off between fitting many voxels at a time, and speeding up the computations. Trying to fit too many will eventually slow the computations, as larger problems take more time to built within the solvers. In our experience, we find that 50 voxels is about the right amount, but this number could exhibit some dependence on the specs of the computer on which computations are performed.

# 3.3 Output: model parameters and scalar invariants

The following call to the software produces the two outputs "model" and "invariants":

```
[model, invariants] = qtiplus_fit(data, btensors);
```

The "model" variable constitutes the main output of the software, and contains the estimated QTI model parameters. These are returned in a 4D Matlab matrix of size [nx, ny, nz, 28] where the arrangement is as follows:

model[x, y, z, 1] =  $S_0$ , the estimated non diffusion-weighted signal, in arbitrary units model[x, y, z, 2:7] = the 6 unique elements of the mean diffusion tensor, in SI units  $[m^2/s]$  model[x, y, z, 8:28] = the 21 unique elements of the covariance tensor, in SI units  $[m^4/s^2]$  with x, y, and z being the indices of a generic voxel in the volume. The functions in the helper functions folder can help in switching between different tensor representations. The second output, "invariants", consists of a structure containing the scalar maps derived from the estimated model parameters. For details about the invariants and their nomenclature please refer to [1]. Figure 2 shows the content of the "invariants" structure. There,  $S_0$  is returned in arbitrary units, while the Mean Diffusivity MD is returned in  $[\mu m^2/ms]$ .

# 3.4 Computational times

An extensive review of the computational times for the different steps of the software and different datasets is provided in [2]. In table 1 we provide exemplary computational times based on the example dataset provided with the software package. The size of the dataset is [64, 64, 6], for a total of 24576 voxels. The machine on which the times were recorded features a 6-core Intel Xeon W-2133 CPU and Matlab R2020b. Computations were performed over 6 workers using the Parallel Computing Toolbox. *nvox* was set to 50.

Table 1: Run times for Mosek and SDPT3				
Solver	Pipeline	Run Times		
Mosek	SDP(dc)	5 min		
Mosek	SDPdc and NLLSdc	11 min		
Mosek	SDPdc, NLLSdc, (m) check, and SDPdcm	14 min		
Mosek	SDPdc, NLLSdc, and SDPdcm	20 min		
Mosek	SDPdc, (m) check, and SDPdcm	8 min		
Mosek	SDPdc and SDPdcm	13 min		
SDPT3	SDPdc	18 min		
SDPT3	SDPdc and NLLSdc	23 min		
SDPT3	SDPdc, NLLSdc, (m) check and SDPdcm	26 min		
SDPT3	SDPdc, NLLSdc, and SDPdcm	41 min		
SDPT3	SDPdc, (m) check, and SDPdcm	20 min		
SDPT3	SDPdc and SDPdcm	35 min		

# 4 Example

In this section we step-by-step go through an example to show how to use the software and the different functionalities. The content of this section is also available in the script example\_qtiplus\_fit.

We start by loading the provided dataset and experimental parameters. The dataset here has size [64, 64, 6, 217].

```
data = niftiread('example_dataset.nii');
load('bten');
```

This dataset is constructed as follows: for the B-tensors contained in the variable "bten", the diffusion MR signal is generated using a non-central Wishart distribution [6] with known mean and covariance. As in [2], two signals are generated, one for a diffusion tensor distribution (DTD) with isotropic mean diffusion tensor, and one for a DTD with anisotropic mean diffusion tensor. Rician distributed signal is then obtained by adding Gaussian noise to the real and imaginary parts of the noiseless signal, creating 4096 signal values at SNR =  $1/\sigma = 30$  and SNR= 15 for both cases, where  $\sigma$  is the standard deviation of the underlying Gaussian distribution.

The data are then organized in a [64, 64, 6] matrix, where each of the 6 slices represents the following:

- 1. Noiseless signals for a non-central Wishart distribution with isotropic mean diffusion tensor
- 2. Signals from 1. with SNR = 30
- 3. Signals from 1. with SNR = 15

- 4. Noiseless signals for a non-central Wishart distribution with anisotropic mean diffusion tensor
- 5. Signals from 4. with SNR = 30
- 6. Signals from 4. with SNR = 15

Note that while this dataset is hardwired, we provide a function called  $create_new_example_dataset$  in the example dataset, with which the user can select the noise level to be applied to slices 2,5 and 3,6. For example, we could create a new dataset where the signals in slices 2 and 5 have SNR = 40, and signals from slices 3 and 6 have SNR = 10. This is done with the following call:

```
data = create_new_example_dataset(bten, 40, 10);
```

Next, we create a mask to be passed to the <a href="qtiplus\_fit">qtiplus\_fit</a> function. While a mask in a real-life case scenario would have the purpose of limiting the computations to the areas of interest, here we use it with the sole purpose of reducing the number of voxels over which computations are performed thus speeding up the fitting on the example dataset. First, we define a variable having the same size as one of the diffusion volumes in our dataset (which in this case is [64, 64, 6]):

```
mask = zeros(64, 64, 6);
```

Then, we select an inner region of size [11, 11] over which computations are to be performed:

```
range = 5;
mask(end/2-range:end/2+range,end/2-range:end/2+range,:) = 1;
```

Next, we decide to exclude some of the collected data from the fitting. This is useful, for example, for excluding outliers present in the data. In our case, we want to exclude the first 13 measurements in the dataset as these were acquired without diffusion weighting. To do so, we define a binary array indicating which of the measurements to keep, and which to discard:

```
indx = ones(217, 1);
indx(1:13) = 0;
```

Now we can call the <a href="qtiplus\_fit">qtiplus\_fit</a> function to perform the fit, and pass the mask and the indices using the respective keywords. Additionally, we specify that we want to perform all the steps in the QTI+ framework, which is done through the keyword *pipeline*, and that we want the fit to be performed over 70 voxels at a time, which is achieved through the keyword *nvox*. Note that since we are not specifying whether we want to use parallel computations, the software will use the default behaviour as explained in section 3.2.5. Moreover, we are also

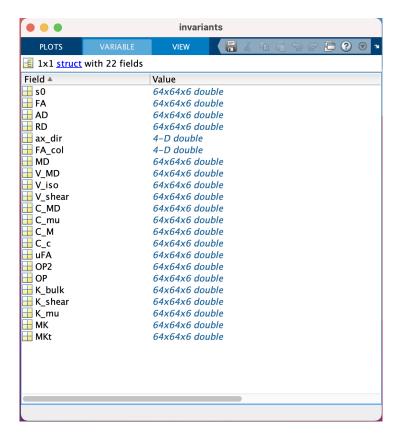


Figure 2: Content of the "invariants" structure as produced by the <a href="qtiplus\_fit">qtiplus\_fit</a> function. For the nomenclature please consult [1].

not specifying which solver to use for the SDP problems, in which case the software will proceed with Mosek if available, and revert to SDPT3 in case it is not. The call to the function will therefore be as follow:

where we assigned the outputs to the variables "model" and "invariants". In Figure 2 the content of the invariants structure is shown. It contains the scalar maps derived from the model parameters.

Using the code provided within the example\_qtiplus\_fit, we can plot one of the scalar maps (the microscopic fractional anisotropy  $\mu$ FA), and look at its value distributions for the different SNRs. This is shown in Figure 3.

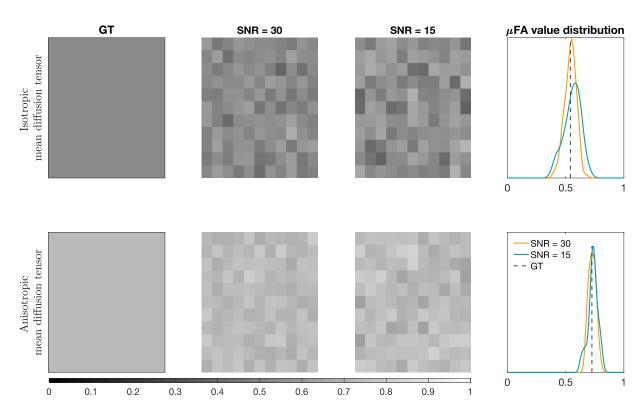


Figure 3: Example maps (here shown is the  $\mu$ FA) obtainable with the software. In the first column, the model is fit to the noiseless signal (slices 1 and 4 in the <code>example\_dataset</code>), thus the produced values can be considered as ground truth. The two middle columns shows the  $\mu$ FA values produced when fitting the model to noisy data at SNR = 30 and SNR = 15, respectively. The rightmost column shows the  $\mu$ FA values distribution for the SNR = 30 and SNR = 15 cases with respect to the ground truth.

## References

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