ANISOTROPY PRESERVING INTERPOLATION OF DIFFUSION TENSORS

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Each voxel of the image contains a symmetric positive definite matrix. The principal characteristics of tensors are [1] - mean diffusivity (characterizes the overall mean squared displacement of molecules) - main direction of diffusivities (orientation in space of the structures of the brain)

degree of anisotropy (describes the degree to which molecular displacements vary in space)

The processing of diffusion tensor images requires the development of methods adapted to 'matrix-valued' images. One of the most used method is the Log-Euclidean framework [2], for which

 $d_{\text{LE}}(\mathbf{S}_1, \mathbf{S}_2) = \parallel \log(\mathbf{S}_1) - \log(\mathbf{S}_2) \parallel$

U,

Proposed similarity measure: based on the spectral decomposition

In this work we propose to use the **spectral decomposition** of diffusion tensors in order to preserve all the diffusion information. Let S_1 and S_2 be two diffusion tensors, whose spectral decompositions are $S_1 = U_1 \Lambda_1 U_1^T$, the proposed similarity measure is

$d_{\rm SD}^{2}(\mathbf{S}_{1},\mathbf{S}_{2}) = k(\Lambda_{1},\Lambda_{2}) d_{\rm SO(3)}^{2}(\mathbf{U}_{1},\mathbf{U}_{2}) + d_{\rm D^{+}(3)}^{2}(\Lambda_{1},\Lambda_{2})$

Spherical or isotropic tensors have low orientational information, so their orientation contribution to the metric should be minimal. $k \in [0,1]$

k is constructed to be small (around zero) when any of the tensor is isotropic and close to 1 when both are highly anisotropic. Distance between orientations is computed through the use of unit quaternions. This parametrization of the rotations enables to speed up the computations.

Because of the non-uniqueness of the spectral decomposition, 4 different rotation matrices represent a single tensor. This implies 8 quaternions for a tensor. The distance between them is computed through

 $d(\mathcal{Q}_1, \mathcal{Q}_2) = \min_{q_2 \in \mathcal{Q}_2} \parallel q_1^r - q_2 \parallel$

This is called the **realignment** of quaternions.

Eigenvalues of tensors are manipulated as **geometric** entities (*i.e.* the logarithms of eigenvalues are used for computations).

The distance between the intensities of diffusion is thus given by

$$d(\Lambda_1,\Lambda_2) = \sqrt{\sum_i \log^2\left(\frac{\lambda_{1,i}}{\lambda_{2,i}}\right)}$$

Mean of a group of N tensors

The mean of a group of N tensors is given by

$$\Lambda_{\mu}$$

The orientation of the mean is given by the mean quaternion.

 $\mathbf{S}_{u} = \mathbf{U}_{u} \boldsymbol{\Lambda}_{u} \mathbf{U}_{u}^{T}$

The eigenvalues of the mean are given by the **geometric mean** of the eigenvalues.

$$\Lambda_{\mu} = \exp(\sum_{i=1}^{N} w_i \log(\Lambda_i))$$

Knowing the mean eigenvalues, the factor k can be computed for each tensor. $k_i = k(\Lambda_u, \Lambda_i)$

Most informative tensor: the one for which the product $w_i k_i$ is the maximum.

$$\mathbf{q}_{m} = \frac{\sum_{i=1}^{w_{i}} w_{i} k_{i} \mathbf{q}_{i,r}}{\sum_{k_{i}} \mathbf{q}_{\mu}} \qquad \mathbf{q}_{\mu} = \frac{\mathbf{q}_{m}}{\|\mathbf{q}_{m}\|}$$

The most informative tensor will be chosen as a reference for the computation of the mean. Each quaternion has to be

This method ensures the **commutativity** and the **uniqueness** of the mean.



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