Computation of electrostatic fields in anisotropic human tissues using the Finite Integration Technique (FIT)

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Abstract. The exposure of human body to electromagnetic fields has in the recent years become a matter of great interest for scientists working in the area of biology and biomedicine. Due to the difficulty of performing measurements, accurate models of the human body, in the form of a computer data set, are used for computations of the fields inside the body by employing numerical methods such as the method used for our calculations, namely the Finite Integration Technique (FIT). A fact that has to be taken into account when computing electromagnetic fields in the human body is that some tissue classes, i.e. cardiac and skeletal muscles, have higher electrical conductivity and permittivity along fibers rather than across them. This property leads to diagonal conductivity and permittivity tensors only when expressing them in a local coordinate system while in a global coordinate system they become full tensors. The Finite Integration Technique (FIT) in its classical form can handle diagonally anisotropic materials quite effectively but it needed an extension for handling fully anisotropic materials. New electric voltages were placed on the grid and a new averaging method of conductivity and permittivity on the grid was found. In this paper, we present results from electrostatic computations performed with the extended version of FIT for fully anisotropic materials.

1 Introduction

tronic devices being used both in households and for communication purposes, special concerns arise regarding the possible adverse biological effects of high or low frequency electromagnetic fields on the human body. The difficulties in directly measuring the fields' induced currents or energy inside the body have resulted in different maximum values within the safety guidelines for limiting the effects of human

With continuously increasing numbers of electrical and elec-

exposure to non-ionizing electromagnetic radiation. In such conditions, the estimation of electromagnetic fields inside the body using numerical methods adequate for computer code implementations, represents a good alternative to the experimental measurements, especially, because the computational capacities are currently able to respond to such complex tasks and continue to have a quick evolution. Another argument for using computer based numerical methods is the availability of realistic human body models of high resolution based on anatomical data in the form of a computer data set. Numerical methods such as the Finite Element Method (FEM), the Finite Difference Method (FD), the Boundary Element Method (BEM), and the Finite Integration Technique (FIT) have already been used in computations of electromagnetic fields in the human body. Our computations are based on an extended version of the Finite Integration Technique (FIT) which allows us to account for the anisotropic character of some tissue classes.

Human body model

The human body model used in our simulations was created by a working group at the Institute of Biomedical Engineering, University of Karlsruhe, Germany, who applied strategies of image processing (Sachse et al., 1996a; 1996b) to a set of anatomical data obtained from Computed Tomography (CT) and Magnetic Resonance Tomography (MRT) provided by the National Library of Medicine, Maryland, USA. The images were first assembled in a 3D anatomical model consisting of cubic voxels each voxel having been assigned one of 40 different tissue types. To every tissue class was assigned a frequency dependent electrical conductivity and permittivity in order to obtain a dielectric model which can be used for electromagnetic simulations in computerized applications. The dielectric properties of tissues are known from measurements (Gabriel et al., 1996). The model, known as HUGO, is available in different resolutions, from 8 mm³ to 1 mm³. Furthermore, this model was extended with the

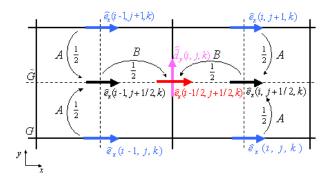


Fig. 1. Local interpolation scheme of a grid voltage in x-direction at the location of a grid flux in y-direction.

Orientation Data Set which gives the direction of fibers by providing two angles for every voxel containing muscle tissue (Sachse et al., 1998).

3 Electromagnetic properties of biological tissues

It is known that biological tissue is non-magnetic i.e. the permeability of biological tissue is equal to that of free space (Durney et al., 1986). The permittivity and conductivity vary with temperature but a stronger variance is experienced with frequency. While permittivity of biological tissue generally decreases with frequency, its conductivity generally increases.

Some tissue types which present a fiber structure (e.g. skeletal and cardiac muscles) are anisotropic having higher conductivity and permittivity in the longitudinal direction of the fibers than on the perpendicular direction to the fibers (Sachse et al., 1997). Shortly, these tissues present a transversely isotropic anisotropy with regard to their dielectric properties.

(Note: In compliance with the subject matter, the remainder of this paper will refer only to permittivity, though, it should be noted that conductivity can be treated in exactly the same way.)

In a local coordinate system, the permittivity of muscle tissue is described by a diagonal tensor of rank two:

$$\mathbf{T}_{\varepsilon_{\mathbf{l}}} = \begin{pmatrix} \varepsilon_{l} & 0 & 0 \\ 0 & \varepsilon_{p} & 0 \\ 0 & 0 & \varepsilon_{p} \end{pmatrix}, \tag{1}$$

where ε_l is the longitudinal permittivity and ε_p is the perpendicular permittivity, relative to the fiber's direction.

Since muscle fibers are miscellaneously oriented in the body, for computational reasons, it is useful to express this tensor in a global coordinate system. In this respect, the local tensor has to be rotated according to the following equation:

$$\mathbf{T}_{\varepsilon_{\mathbf{G}}} = \mathbf{R} \mathbf{T}_{\varepsilon_{\mathbf{L}}} \mathbf{R}^{-1}, \tag{2}$$

where the rotation matrix **R** is the product of two other rotation matrices: $\mathbf{R} = \mathbf{R}_{xy} \mathbf{R}_{xz}$, with \mathbf{R}_{xy} and \mathbf{R}_{xz} given below.

$$\mathbf{R}_{xy} = \begin{pmatrix} \cos \phi - \sin \phi & 0 \\ \sin \phi & \cos \phi & 0 \\ 0 & 0 & 1 \end{pmatrix} \quad \mathbf{R}_{xz} = \begin{pmatrix} \sin \theta & 0 & \cos \theta \\ 0 & 1 & 0 \\ \cos \theta & 0 - \sin \theta \end{pmatrix}. \quad (3)$$

The angles ϕ and θ are the rotation angles about the z- and y-axes, respectively, provided by the Orientation Data Set. After the rotation in Eq. (2) we obtain a full and symmetric tensor which expresses the permittivity of muscle tissue in a global coordinate system.

4 Considerations concerning the classical Finite Integration Technique (FIT)

Since a lot of literature has already been published about the FIT, only a short introduction is provided here concerning the allocation of some grid-state variables and material treatment related to the subject of this paper. For more general details about FIT, we recommend you see van Rienen (2001), meanwhile, for more information concerning the subject matter of this paper, see van Rienen et al. (2003).

The Finite Integration Technique (FIT) was first published by Weiland (1977) and developed as a numerical method which discretizes the Maxwell's equations on a grid pair preserving the analytical properties of the original equations (van Rienen, 2001). On the FIT's grid doublet, in connection with Electrostatics, the following variables are defined:

- the electric potentials (denoted with φ_n), allocated in every mesh node belonging to the primary grid;
- the electric grid voltages (denoted with \hat{e}_n), allocated in the middle of the edges belonging to the primary grid and calculated as the integral of the electric field along primary edges;
- the electric grid flux densities (denoted with \hat{d}_n), normal in the middle of the surfaces belonging to the dual grid and calculated as the integral of the electric flux on dual surfaces.

In a 3D Cartesian coordinate system, because the dual grid is shifted with half an edge length in all positive directions with respect to the primary grid, the primary edges intersect dual surfaces on the normal direction in the middle. This means that an electric grid voltage \hat{e}_n from the primary grid is allocated in the same point with the corresponding electric grid flux \hat{d}_n in the same direction, from the dual grid.

The electric fluxes and voltages with coinciding both, locations and orientations, are related to each other by the permittivity according to the following equation:

$$\frac{\hat{d}_{n}}{\hat{e}_{n}} = \frac{\int \int_{\tilde{A}_{n}} \mathbf{D} \cdot d\mathbf{A}}{\int_{L_{n}} \mathbf{E} \cdot d\mathbf{s}} = \frac{\int \int_{\tilde{A}_{n}} \varepsilon d\mathbf{A} + O(\Delta^{\kappa+1})}{\int_{L_{n}} d\mathbf{s} + O(\Delta^{\kappa})}$$

$$\approx \frac{\overline{\varepsilon} \int \int_{\tilde{A}_{n}} d\mathbf{A}}{\int_{L_{n}} d\mathbf{s}} + O(\Delta^{\kappa}) \approx \overline{\varepsilon} \frac{|\tilde{A}_{n}|}{|L_{n}|} = [M_{\varepsilon}]_{n,m}, \tag{4}$$

where κ takes values between $\kappa=2$ for varying permittivity or non-uniform step size and $\kappa=3$ otherwise. The symbol $\overline{\epsilon}$ denotes a weighted average of the permittivity on a dual surface from four possibly different values belonging to the primary grid cells intersected by that dual surface. For the entire grid, the point-wise relation in Eq. (4) becomes:

$$\hat{\hat{d}} = \tilde{D}_{\mathbf{A}} \mathbf{D}_{\varepsilon} \mathbf{D}_{\mathbf{S}}^{-1} \hat{\boldsymbol{e}} = \mathbf{M}_{\varepsilon} \hat{\boldsymbol{e}}, \tag{5}$$

where: $\tilde{\mathbf{D}}_{\mathbf{A}}$, \mathbf{D}_{ε} and $\mathbf{D}_{\mathbf{S}}^{-1}$ are diagonal matrices containing the areas of the dual surfaces, the permittivity of the primary grid cells averaged on the dual surfaces and the inverse of the primary edge lengths, respectively. The vectors $\hat{\boldsymbol{d}}$ and $\hat{\boldsymbol{e}}$ contain the electric grid fluxes allocated in the middle of dual surfaces and the electric grid voltages along the primary edges. \mathbf{M}_{ε} is the material operator which decomposed along the axes of a Cartesian coordinate system has the following diagonal form:

$$\mathbf{M}_{\varepsilon} = \begin{pmatrix} \tilde{\mathbf{D}}_{\mathbf{A}_{yz}} \ \mathbf{D}_{\varepsilon_{xx}} \ \mathbf{D}_{\mathbf{S}_{x}}^{-1} & \mathbf{0} & \mathbf{0} \\ \mathbf{0} & \tilde{\mathbf{D}}_{\mathbf{A}_{xz}} \ \mathbf{D}_{\varepsilon_{yy}} \ \mathbf{D}_{\mathbf{S}_{y}}^{-1} & \mathbf{0} \\ \mathbf{0} & \mathbf{0} & \tilde{\mathbf{D}}_{\mathbf{A}_{xy}} \ \mathbf{D}_{\varepsilon_{zz}} \ \mathbf{D}_{\mathbf{S}_{z}}^{-1} \end{pmatrix}$$
(6)

In its classical form, FIT allows the presence of diagonally anisotropic material on the grid but, as it was shown, this allowance is not enough for computing muscle tissues which are fully anisotropic in a global coordinate system.

5 Extension of the Finite Integration Technique for computing anisotropic tissues

To deal with anisotropic tissues, we follow the idea from (Krüger, 2000) where gyrotropic materials were treated in time domain.

When the diagonal matrix \mathbf{D}_{ε} in Eq. (5) is replaced with a full one, the material operator in Eq. (6) becomes:

$$\mathbf{M}_{\varepsilon} = \begin{pmatrix} \tilde{\mathbf{D}}_{\mathbf{A}_{yz}} \ \mathbf{D}_{\varepsilon_{xx}} \ \mathbf{D}_{\varepsilon_{x}}^{-1} & \tilde{\mathbf{D}}_{\mathbf{A}_{yz}} \ \mathbf{D}_{\varepsilon_{xy}} \ \mathbf{D}_{\varepsilon_{x}}^{-1} & \tilde{\mathbf{D}}_{\mathbf{A}_{yz}} \ \mathbf{D}_{\varepsilon_{xz}} \ \mathbf{D}_{\varepsilon_{xz}}^{-1} \\ \tilde{\mathbf{D}}_{\mathbf{A}_{xz}} \ \mathbf{D}_{\varepsilon_{yx}} \ \mathbf{D}_{S_{x}}^{-1} & \tilde{\mathbf{D}}_{\mathbf{A}_{xz}} \ \mathbf{D}_{\varepsilon_{yy}} \ \mathbf{D}_{S_{y}}^{-1} & \tilde{\mathbf{D}}_{\mathbf{A}_{xz}} \ \mathbf{D}_{\varepsilon_{yz}} \ \mathbf{D}_{S_{z}}^{-1} \\ \tilde{\mathbf{D}}_{\mathbf{A}_{xy}} \ \mathbf{D}_{\varepsilon_{xz}} \ \mathbf{D}_{S_{x}}^{-1} & \tilde{\mathbf{D}}_{\mathbf{A}_{xy}} \ \mathbf{D}_{\varepsilon_{zz}} \ \mathbf{D}_{S_{z}}^{-1} \end{pmatrix}$$
(7)

The off-diagonal terms of the material operator are coupling electric grid flux vectors in one direction to electric grid voltage vectors in another direction. To bring the vectors with different orientations to the same location on the grid, an interpolation process is necessary (Krüger, 2000) for which a local scheme is presented in Fig. 1.

In Fig. 1, the electric grid voltage $\hat{e}_x(i-1/2, j+1/2, k)$ is interpolated in two steps at the location of $\hat{d}_y(i, j, k)$. In the first step, every pair of electric voltages having the same coordinate on the x-axis, are interpolated along the y-axis in the middle (each voltage contributing with a factor of a half) through an A-type interpolation, building the voltages $\hat{e}_x(i-1, j+1/2, k)$ and $\hat{e}_x(i, j+1/2, k)$. In the second step, these two voltages are interpolated along the x-axis in the middle (each contributing with a factor of a

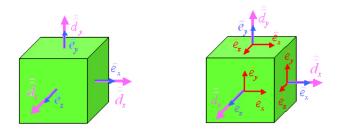


Fig. 2. Allocation of both electric grid voltages and electric grid fluxes relative to a dual FIT cell. Left: before the interpolation. Right: after the interpolation.

half) through a B-type interpolation, building the voltage $\hat{e}_x(i-1/2, j+1/2, k)$.

The local interpolation process is expressed by the following equation:

$$\hat{e}_x(i-1/2, j+1/2, k) = \frac{1}{4} [\hat{e}_x(i, j, k) + \hat{e}_x(i, j+1, k) + \hat{e}_x(i-1, k, k) + \hat{e}_x(i-1, j+1, k)].$$
(8)

Figure 2 shows the allocation of voltages relative to a dual cell after the interpolation (right-hand side) compared to the classical allocation (left-hand side).

To globally account the interpolation process, the following matrices are defined:

$$[Q_x]_{pq} = \begin{cases} \frac{1}{2} \ p = q \text{ or } p = q + 1\\ 0 \text{ else} \end{cases}$$
 (9)

$$\left[Q_{y}\right]_{pq} = \begin{cases} \frac{1}{2} & p=q \text{ or } p=q+I\\ 0 & \text{else} \end{cases}$$
 (10)

$$[Q_z]_{pq} = \begin{cases} \frac{1}{2} & p = q \text{ or } p = q + IJ \\ 0 & \text{else} \end{cases} , \tag{11}$$

where I and J represent the maximum number of grid nodes in x- and y-directions, respectively (Krüger, 2000). Within the FIT algorithm, the interpolation matrices will be placed at the off-diagonal terms in the material operator.

In order to keep the symmetry of the physical permittivity tensor the off-diagonal terms $[\mathbf{M}_{\varepsilon}]_{\mathbf{u},\mathbf{v}}$ and $[\mathbf{M}_{\varepsilon}]_{\mathbf{v},\mathbf{u}}$ of the discrete material operator have to be equal. The symmetry of the material operator can be reached through the following steps:

Assuming that primary and dual edges in the same direction have the same length, their cancellation leads to equal grid information in the above-mentioned terms of the material operator according to Eq. (12).

$$\frac{\tilde{A}_{vw}}{L_v} = \frac{\tilde{L}_v \tilde{L}_w}{L_v} = \frac{\tilde{A}_{uw}}{L_u} = \frac{\tilde{L}_u \tilde{L}_w}{L_u} \approx \tilde{L}_w. \tag{12}$$

The approximation introduced in Eq. (12) is given by $L_u = \tilde{L}_u + O(\Delta)$, where the first order error term vanishes if the grid is equidistant. This condition can be applied when meshing the human body model because it consists of cubic voxels.

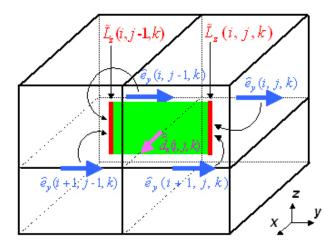


Fig. 3. Construction of electric voltages on the grid.

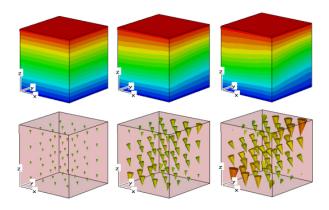


Fig. 4. Top row: electric potential distribtuion. Bottom row: electric flux density. Left column: $\varepsilon_l/\varepsilon_p=1$. Middle column: $\varepsilon_l/\varepsilon_p=2$. Right column: $\varepsilon_l/\varepsilon_p=3$

- Averaging the permittivity along the common dual edge \tilde{L}_w resulting from the first step leads to equal material information in the considered terms. The permittivity average for each off-diagonal term of the material operator uses the approximation from Eq. (12):

$$\frac{\iint_{\tilde{A}_{uw}} \varepsilon \, dA + O(\Delta^{k+1})}{\int_{L_u} ds + O(\Delta^k)} = \frac{\bar{\varepsilon} \left| \tilde{L}_w \right| \int_{\tilde{L}_u} ds}{\int_{L_u} ds} + O(\Delta^k)
\approx \bar{\varepsilon} \left| \tilde{L}_w \right| + O(\Delta^{k^*}) \approx \bar{\varepsilon} \left| \tilde{L}_w \right|,$$
(13)

where $k^* \in [1, 3]$.

Positioning the interpolation matrices within the material operator such as to ensure the symmetry regarding the interpolation process.

Table 1. Electric flux density (max value [C/m²]) given in % relative to the max value in the case of $\varepsilon_I/\varepsilon_p$ =3.

		$\varepsilon_l/\varepsilon_p=1$	$\varepsilon_l/\varepsilon_p=2$	$\varepsilon_l/\varepsilon_p=3$
6	$l_{\text{max}} [\text{C/m}^2]$	13.5%	60.8%	100%

Taking into account the above considerations, the material operator becomes:

$$\mathbf{M}_{\varepsilon} = \begin{pmatrix} \tilde{\mathbf{D}}_{\mathbf{A}_{yz}} \ \mathbf{D}_{\varepsilon_{xx}} \ \mathbf{D}_{\varepsilon_{x}}^{-1} & \mathbf{Q}_{\mathbf{S}}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{xy}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{z}} \ \mathbf{Q}_{x} & \mathbf{Q}_{z}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{xx}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{y}} \ \mathbf{Q}_{x} \\ \mathbf{Q}_{x}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{yx}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{z}} \ \mathbf{Q}_{y} & \tilde{\mathbf{D}}_{\mathbf{A}_{xz}} \ \mathbf{D}_{\varepsilon_{yy}} \ \mathbf{D}_{\varepsilon_{y}}^{-1} & \mathbf{Q}_{z}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{yz}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{x}} \ \mathbf{Q}_{y} \\ \mathbf{Q}_{x}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{zx}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{y}} \ \mathbf{Q}_{z} & \mathbf{Q}_{z}^{\mathrm{T}} \ \mathbf{D}_{\varepsilon_{zx}} \ \tilde{\mathbf{D}}_{\mathbf{S}_{x}} \ \mathbf{Q}_{z} & \tilde{\mathbf{D}}_{\mathbf{A}_{xy}} \ \mathbf{D}_{\varepsilon_{zz}} \ \mathbf{D}_{\mathbf{S}_{z}}^{-1} \end{pmatrix} . (14)$$

This distribution of the interpolation matrices to the off-diagonal terms ensures the global symmetry of the material operator. Due to the permittivity average, when coupling two directions the A-type interpolation is weighted with the length of a dual edge on the third direction. To keep the symmetry, the B-type interpolation cannot be length-weighted. Figure 3 shows that the voltages $\hat{e}_y(i, j-1, k)$ and $\hat{e}_y(i+1, j-1, k)$ are built along the dual edge $\tilde{L}_z(i, j-1, k)$ and the voltages $\hat{e}_y(i, j, k)$, $\hat{e}_y(i+1, j, k)$ are built along the dual edge $\tilde{L}_z(i, j, k)$.

6 Simulations

The human body model (HUGO) offered by the simulation software package CST EMStudioTM (CST GmbH) was used for the import of a cubic volume of muscle tissues within a C++ code where the anisotropic FIT was implemented. The diagonal direction of muscle fibres is described by the angles $\phi=\theta=45^{\circ}$. These muscles were placed in the electrostatic field determined by the imposed potential values of $-10 \, \text{V}$ and $+10 \, \text{V}$, at the boundary planes $Z=\min$ and $Z=\max$, through the Dirichlet boundary conditions. All the other boundaries, were treated with Neumann boundary conditions. Three simulations were performed using the CG solver from PETSc (Balay et al., 1997) corresponding to three different values for the ratio between longitudinal and perpendicular permittivity of muscle tissues with respect to their fibre direction, i.e. $\varepsilon_l/\varepsilon_p=1$; 2; 3.

7 Results and discussions

In Fig. 4, the top row presents three scalar plots of the electric potentials computed inside a muscle volume considered to be isotropic (left), anisotropic (having the ratio $\varepsilon_l/\varepsilon_p$ =2 (middle)) and anisotropic (with $\varepsilon_l/\varepsilon_p$ =3 (right)). Comparing the plots, the equipotential planes were aligned to the direction of muscle fibers (diagonal direction with regard to the Cartesian coordinate system) in the anisotropic cases. The magnitude of this alignment depends on the magnitude of the ratio between longitudinal and perpendicular permittivity ($\varepsilon_l/\varepsilon_p$).

The bottom row in Fig. 4 presents three vector plots of the electric flux density, each computed with the corresponding electric potentials in the same column and scaled to the same maximum value which was found in the anisotropic case of $\varepsilon_l/\varepsilon_p$ =3. Relative to this value, the other maximum values are given in percentage in Table 1. Compared to the isotropic case where the electric flux is homogeneous in the entire muscle volume, in the anisotropic cases, the flux has higher values along the fiber's direction.

8 Conclusions

In this paper we presented the anisotropy of muscle tissues with regard to dielectric permittivity and its mathematical analytic model. This model was further discretized to conform with the numerical algorithm of the Finite Integration Technique and implemented in software code. With this code, more simulations were performed and the results were compared to the isotropic case.

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