

# Comparison of Two Formulations for Computing Body Surface Potential Maps

Emma Lagracie, Lisl Weynans, Yves Coudière

Univ. Bordeaux, CNRS, Bordeaux INP, IMB, UMR 5251, IHU Liryc  
F-33400 Talence, France

We want to compute QRS complexes from activation maps and a predefined shape for the transmembrane voltage  $v$ . In the context of ECGi, it would allow to directly reconstruct activation maps, without resorting to reconstruction and post-processing of cardiac electrical potentials. Here we compare two possible formulations to compute QRS complexes.

Several methods exist to simulate an ECG. Assuming that the transmembrane voltage  $v$  is given in the heart  $H$ , the electrical potential in the torso (extracardiac  $T$  and cardiac extracellular  $H$ )  $u$  may be retrieved following two approaches. Either it solves a Laplace equation with discontinuous conductivity coefficient (heart and torso) and ionic current as a source (source formulation (1)), specifically,

$$\begin{aligned} -\operatorname{div}(\sigma_e \nabla u) &= \partial_t v + f(v) && \text{in } H, \\ -\operatorname{div}(\sigma_T \nabla u) &= 0 && \text{in } T; \end{aligned} \quad (1)$$

or the quasi-stationary electrical balance between the intra and extracellular fields (balance formulation (2)), specifically

$$\begin{aligned} -\operatorname{div}(\sigma_e \nabla u) - \operatorname{div}(\sigma_i \nabla(u + v)) &= 0 && \text{in } H, \\ -\operatorname{div}(\sigma_T \nabla u) &= 0 && \text{in } T. \end{aligned} \quad (2)$$

Commonly, the potential  $u$  is computed from the balance formulation. Anyway, we may also use the source formulation. Note that both formulations coincide only if  $v$  solves the complete bidomain equations. We compute reference activation maps and QRS complexes with a bidomain code. Afterwards the electrical field  $u$  is also calculated from the activation maps and a predefined  $v$  (and ionic model  $f(v)$ ), using the two different formulations.

We compare the potential fields  $u$  and the QRS complexes obtained by these two methods to the reference ones. Preliminary results show some significant differences between the two methods, with a better accuracy for the most popular balance formulation (2), for a smoothed heaviside form of  $v$ .