

Supplementary Materials

Irreversible Inactivation of Initial Form of Water-Soluble Redox Proteins-Theoretical Study in Square-Wave Voltammetry

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$$E_s := 0.4 \quad \Delta E := 0.8 \quad dE := 0.004 \quad E_{sw} := 0.05$$

$$n := 1 \quad \frac{F}{RT} := 96500 \quad R_{sw} := 8.314 \quad T_{sw} := 298.15 \quad D := 0.000005$$

$$j := 1, \frac{\Delta E}{dE} \cdot 50 \quad \alpha := 0.5$$

$$k_c := 30055.2500$$

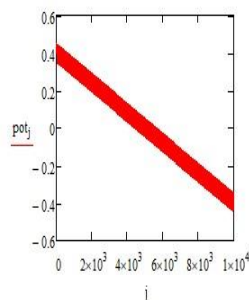
$$f := 10$$

$$k_s := 10^{-1.50504499512}$$

Model of Diffusional Electrode Mechanism with Irreversible Chemical Reaction Coupled to Initial Redox Form in Square-Wave Voltammetry

$$pot_j := E_s + E_{sw} - \left[\text{cell} \left(\frac{j}{25}, \frac{1}{2} \right) \cdot dE + \text{if} \left(\frac{\text{cell} \left(\frac{j}{25} \right)}{2} = \text{cell} \left(\frac{j}{25}, \frac{1}{2} \right) \cdot 1, -1 \right) \cdot E_{sw} + E_{sw} \right] - dE$$

$$K_{et} = \frac{k_s}{f^{0.5} \cdot D^{0.5}} \quad K_{et} = 100.002$$



$$k := 1, \frac{\Delta E}{dE} \cdot 50$$

$$K_{chem} = \frac{k_c}{f} \quad K_{chem} = 3.006 \times 10^3$$

$$\Phi_{\text{red}} := n \cdot \frac{F}{R \cdot T} \cdot pot_j$$

$$S_{\text{sk}} := \sqrt{k} - \sqrt{k-1}$$

$$\frac{I_{\text{sk}}}{I_{\text{sk}}} := \text{erfc} \left[\left[K_{chem} \cdot \frac{k}{50} \right]^{0.5} \right] - \text{erfc} \left[\left[K_{chem} \cdot \frac{(k-1)}{50} \right]^{0.5} \right]$$

Definitions and Meanings of the Symbols

- f** is the SW frequency
- k_s** is standard rate constant of electron transfer
- α** is electron transfer coefficient
- n** is number of exchanged electrons
- dE** is potential step
- E_{sw}** is square-wave amplitude
- T** is thermodynamic temperature
- R** is universal gas constant
- k_c** is rate constant of irreversible chemical reaction
- K_{et}** is dimensionless kinetic electrode parameter
- K_{chem}** is dimensionless kinetic chemical parameter
- S_k** is numerical integration factor
- E_s** is starting potential
- Φ** is dimensionless potentials
- F** is the Faraday constant
- Ψ** is dimensionless current

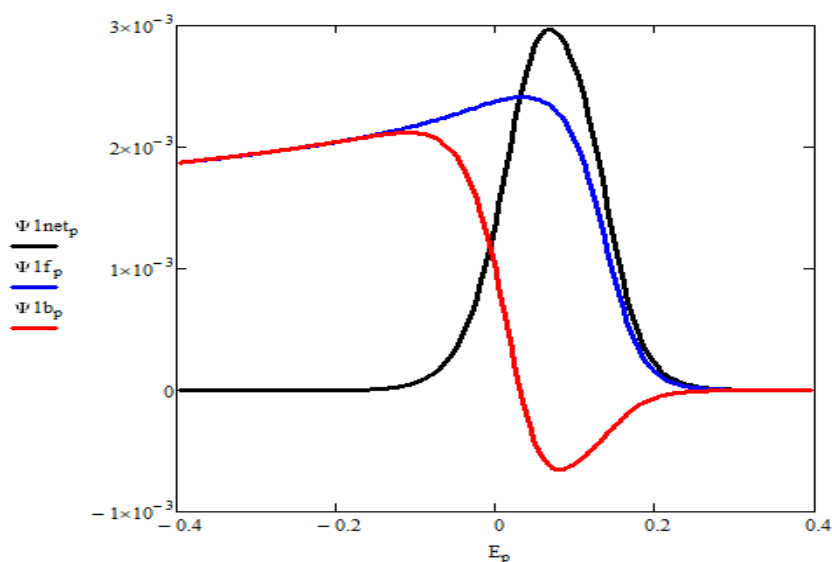
$$\Psi_{1_1} := \frac{\frac{\text{Ket} \cdot e^{-\alpha \cdot \Phi_1}}{(50 \cdot \pi)^{0.5}} - \left[\frac{2\text{Ket} \cdot e^{-\alpha \cdot \Phi_1} \cdot (1 + e^{\Phi_1})}{(50 \cdot \pi)^{0.5}} \cdot 0 + \frac{\text{Kchem}^1 \cdot e^{-\alpha \cdot \Phi_1} \cdot 1 \cdot L_1}{(50 \cdot \pi)^{0.5}} \right]}{1 + \frac{2\text{Ket} \cdot e^{-\alpha \cdot \Phi_1} \cdot (1 + e^{\Phi_1}) \cdot S_1}{(50 \cdot \pi)^{0.5}} - \frac{\text{Kchem}^1 \cdot e^{-\alpha \cdot \Phi_1} \cdot 1 \cdot L_1}{(50 \cdot \pi)^{0.5}}}$$

$$\Psi_{1_k} := \frac{\frac{\text{Ket} \cdot e^{-\alpha \cdot \Phi_k}}{(50 \cdot \pi)^{0.5}} + \frac{\text{Kchem}^1 \cdot e^{-\alpha \cdot \Phi_k} \cdot 1}{(1)^{0.5}} \cdot \sum_{j=1}^{k-1} (\Psi_{1_j} \cdot L_{k-j+1}) - 2\text{Ket} \cdot e^{-\alpha \cdot \Phi_k} \cdot \frac{(1 + e^{\Phi_k})}{(50 \cdot \pi)^{0.5}} \cdot \sum_{j=1}^{k-1} (\Psi_{1_j} \cdot S_{k-j+1})}{1 + \frac{2\text{Ket} \cdot e^{-\alpha \cdot \Phi_k} \cdot (1 + e^{\Phi_k}) \cdot S_1}{(50 \cdot \pi)^{0.5}} - \frac{\text{Kchem}^1 \cdot e^{-\alpha \cdot \Phi_k} \cdot 1 \cdot L_1}{(1)^{0.5}}}$$

$$p := 1 \cdot \left(\frac{\Delta E}{dE} \right) - 1$$

$$\Psi_{1f_p} := \Psi_{1_{(p+1) \cdot 50}} \quad \Psi_{1b_p} := \Psi_{1_{50 \cdot p + 25}} \quad \Psi_{1net_p} := \Psi_{1f_p} - \Psi_{1b_p}$$

$$E_p := E_s - p \cdot dE$$



References:

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4. R. Gulaboski, E. S. Ferreira, C. M. Pereira, M. N. D. S. Cordeiro, A. Garau, V. Lippolis, A. F. Silva, *Journal of Physical Chemistry C*, 112 (2008) 153-161