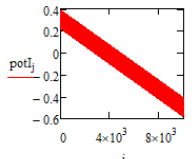


**Supplementary Material: Electrochemically induced dimerization of lipophilic redox proteins: theoretical insights in protein-film square-wave voltammetry: *Monatsh Chem* (2023). <https://doi.org/10.1007/s00706-023-03065-4>**

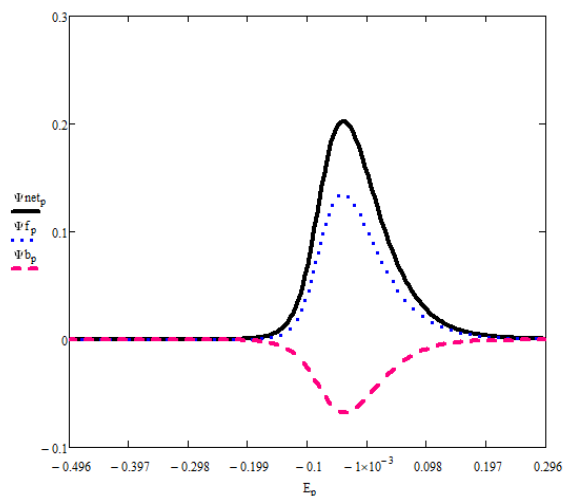
**Sanja Lazarova, Pavle Apostoloski & Rubin Gulaboski**

$E_{sl} := 0.3$     $\Delta E := 0.8$     $dE := 0.004$     $E_{sw} := 0.085$     $ks := 2.5$     $kf := 0.0005$     $cOx := 10.0$     $rs := 1$   
 $n := 1$     $F := 96500$     $R := 8.314$     $T := 298.15$     $\alpha := 0.5$     $f := 10$     $kb := 0.0005$     $cs := \frac{cOx}{rs}$   
 $j := 1.. \frac{\Delta E}{dE} \cdot 50$     $Ket := \frac{ks}{f}$     $Keq := 1.00$    **MATHCAD FILE**  
 $pot_j := E_{sl} + E_{sw} - \left[ \left[ \text{ceil} \left( \frac{j}{25} \cdot \frac{1}{2} \right) dE + \text{if} \left( \frac{\text{ceil} \left( \frac{j}{25} \right)}{2} = \text{ceil} \left( \frac{j}{25} \cdot \frac{1}{2} \right), 1, -1 \right) \cdot E_{sw} + E_{sw} \right] - dE \right]$   
 $K_{chem} := \frac{(kf + kb)}{f}$   
 $M_j := e^{-K_{chem} \cdot \frac{j}{50}} - e^{-K_{chem} \cdot \frac{j+1}{50}}$   
  
 $\Phi_{I_j} := n \cdot \frac{F}{R \cdot T} \cdot pot_j$   
 $\Psi_{I_1} := \frac{Ket \cdot e^{-\alpha \cdot \Phi_{I_1}}}{1 + \frac{Ket}{50} \cdot e^{-\alpha \cdot \Phi_{I_1}} + \frac{1}{cs} \cdot Ket \cdot K_{chem}^{-1} \cdot e^{-\Phi_{I_1} \cdot (1-\alpha)} \cdot M_1}$   
 $\Psi_{I_j} := \frac{\frac{Ket}{cs} \cdot e^{-\alpha \cdot \Phi_{I_j}} - \frac{Ket}{50 \cdot cs} \cdot e^{-\alpha \cdot \Phi_{I_j}} \cdot \sum_{i=1}^{j-1} \Psi_{I_i} - \left( \frac{1}{1} \right) \left[ 1 - \frac{Ket}{1} \cdot K_{chem}^{-1} \cdot e^{-\Phi_{I_j} \cdot (1-\alpha)} \cdot \sum_{i=1}^{j-1} (\Psi_{I_i} \cdot M_j) - \frac{1 \cdot Keq}{1 + Keq} \cdot K_{chem}^{-1} \cdot e^{-\Phi_{I_j} \cdot (1-\alpha)} \cdot \sum_{i=1}^{j-1} (\Psi_{I_i} \cdot M_j) \right]^2}{1 + \frac{Ket}{50 \cdot cs} \cdot e^{-\alpha \cdot \Phi_{I_j}} + \frac{1}{1} \left[ 1 - \frac{Ket}{1} \cdot K_{chem}^{-1} \cdot e^{-\Phi_{I_j} \cdot (1-\alpha)} \cdot M_1 - \frac{1 \cdot Keq}{1 + Keq} \cdot K_{chem}^{-1} \cdot e^{-\Phi_{I_j} \cdot (1-\alpha)} \cdot M_1 \right]^2}$   
 $\Psi_j := \Psi_{I_j}$     $p := 1.. \left( \frac{\Delta E}{dE} \right) - 1$   
 $\Psi_{b_p} := \Psi_{50 \cdot p + 25}$     $\Psi_{f_p} := \Psi_{(p+1) \cdot 50}$     $\Psi_{net_p} := \Psi_{f_p} - \Psi_{b_p}$     $E_p := E_{sl} - p \cdot dE$

Surface E-dimerization mechanism in SWV  
 $Ox(ads) + 1e^- = Red(ads)$   
 $2Red(ads) \rightleftharpoons D$

Meaning of symbols in the MATHCAD File

- ks** is standard rate constant of electron transfer (s<sup>-1</sup>)
- kf** is forward rate constant of dimerization step
- kb** is reverse rate constant of dimerization step
- f** is the SW frequency (Hz)
- Esw** is SW amplitude (V)
- dE** is potential step (V)
- n** is number of electrons exchanged
- α** is electron transfer coefficient
- F** is Farady constant
- R** is universal gas constant
- T** is thermodynamic temperature
- cOx** is molar concentration of Ox
- rs** is auxiliary constant
- cs** is dimensionless concentration parameter
- Kchem** is chemical rate parameter of dimerization reaction
- Keq** is equilibrium constant of dimerization reaction
- Mj** is numerical integration factor
- ΦIj** is dimensionless potential
- Esl** is starting potential
- Ket** is dimensionless parameter related to electron transfer step
- Ψ** is symbol of the dimensionless current



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