## **Supplementary Information to**

## Aβ Self-Association and Adsorption on a Hydrophobic

## Nanosurface: Competitive Effects and the Detection of Small

## **Oligomers via Electrical Response**

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Figure S1. Time evolution of the volume of the simulated box  $(Å^3)$  in surface adsorbed a) monomeric, b) dimeric, c) trimeric systems during equilibration in the presence of KCl ions, prior to the application of electric field.



Figure S2. Rotational time correlation functions,  $\langle \mathbf{R}(0) \cdot \mathbf{R}(t) \rangle$ , of the unit vector  $\mathbf{R}(t)$ , joining the  $C_{\alpha}$  atoms of the first and the last residues of the A $\beta$  units, in the monomeric, dimeric and trimeric systems. Data for the free systems are shown in black; for the SWCNT surface adsorbed systems in orange; and for the surface adsorbed systems in the presence of a potential difference of 2.25 Volts in purple. For the dimeric and trimeric systems, the time correlation functions have been averaged over the multiple peptide units.