



Mechanistic Insights into the Release of Doxorubicin from Graphene Oxide in Cancer Cells

Erica Quagliarini ^{1,+}, Riccardo Di Santo ^{2,+}, Daniela Pozzi ^{2,*}, Paolo Tentori ^{3,4}, Francesco Cardarelli ⁴ and Giulio Caracciolo ^{2,*}

- ¹ Department of Chemistry, Sapienza University of Rome, P.le A. Moro 5, 00185 Rome, Italy; erica.quagliarini@uniroma1.it
- ² Department of Molecular Medicine, Sapienza University of Rome, Viale Regina Elena 291, 00161 Rome, Italy; riccardo.disanto@uniroma1.it
- ³ Center for Nanotechnology Innovation@NEST (CNI@NEST), Istituto Italiano di Tecnologia, Piazza San Silvestro 12, 56127 Pisa, Italy; paolo.tentori@sns.it
- ⁴ NEST Laboratory, Scuola Normale Superiore, Piazza San Silvestro 12, 56127 Pisa, Italy; francesco.cardarelli@sns.it
- * Correspondence: giulio.caracciolo@uniroma1.it (G.C); daniela.pozzi@uniroma1.it (D.P).
- + These authors contributed equally to this work.



Figure S1. (a) Absorbance spectra of doxorubicin (DOX) at several known concentrations. (b) DOX calibration curve was generated from absorbance values at λ = 480 nm. The calibration curve was obtained by plotting the absorbance 480 nm versus DOX concentration and was treated by linear regression analysis (dashed line).





Figure S2. Hydrodinamic diameter (DH) of GO/DOX complexes as a function of pH solution.



Figure S3. DH intensity-weighted distribution of pristine GO and GO/DOX complexes.



Figure S4. (a) The red circle highlights the experimental lifetime of DOX in aqueous solution (1 ns). The blue circle highlights the experimental lifetime of pristine graphene-oxide (GO) in aqueous solution. (b) The experimental cluster of GO-DOX is a linear combination of the two components, GO and DOX (represented by the red and blue dots).