

# **One-pot carboxyl enrichment fosters water-dispersibility of reduced graphene oxide: a combined experimental and theoretical assessment**

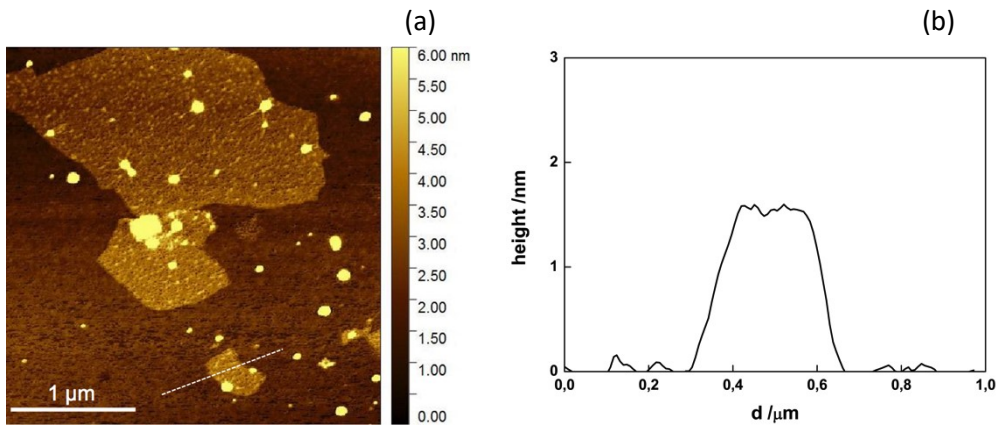
Francesco Amato,\*<sup>1</sup> Alessandro Motta,<sup>1,2</sup> Leonardo Giaccari,<sup>1</sup> Roberto Di Pasquale,<sup>1</sup> Francesca Anna Scaramuzzo,<sup>3</sup> Robertino Zanoni,<sup>1</sup> Andrea Giacomo Marrani\*<sup>1</sup>

<sup>1</sup>*Dipartimento di Chimica, Università di Roma La Sapienza, p.le A. Moro 5, I-00185 Rome, Italy*

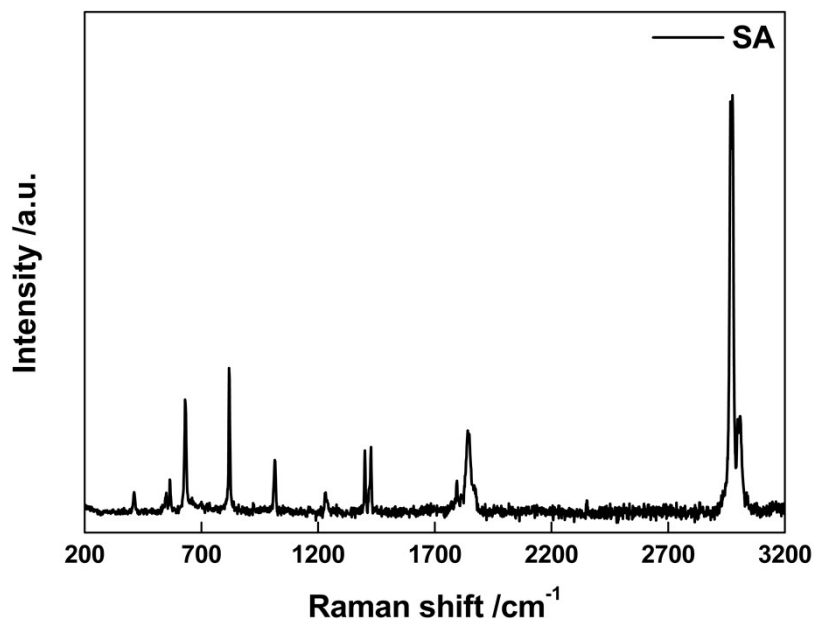
<sup>2</sup>*Consorzio INSTM UdR Roma “La Sapienza”, p.le A. Moro 5, I-00185 Rome, Italy*

<sup>3</sup>*Dipartimento di Scienze di Base e Applicate per l'Ingegneria (S.B.A.I.), Università di Roma La Sapienza, Via del Castro Laurenziano 7, I-00161 Rome, Italy*

## **SUPPLEMENTARY MATERIAL**



**Figure S1.** (a) AFM-tapping mode image of GO. (b) Height profile of the dashed line of figure a.

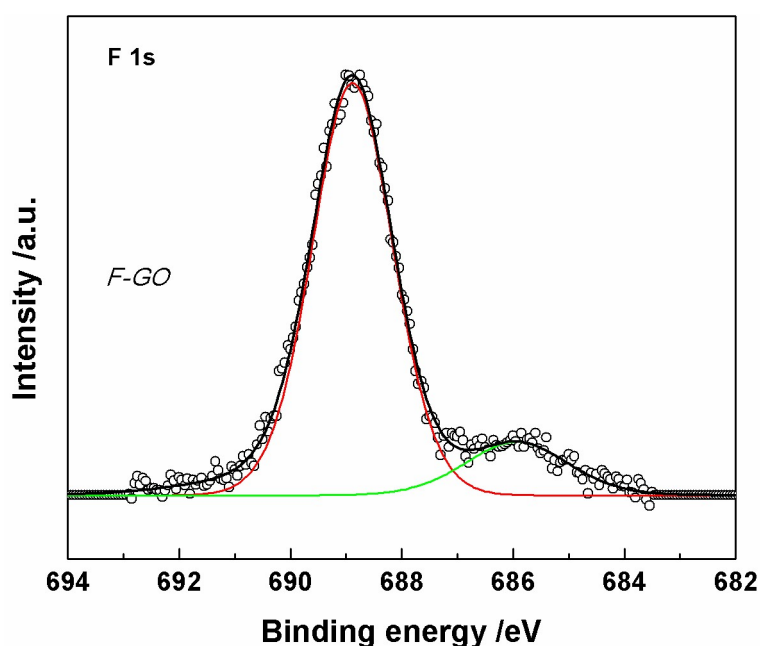


**Figure S2.** Raman spectrum of succinic anhydride

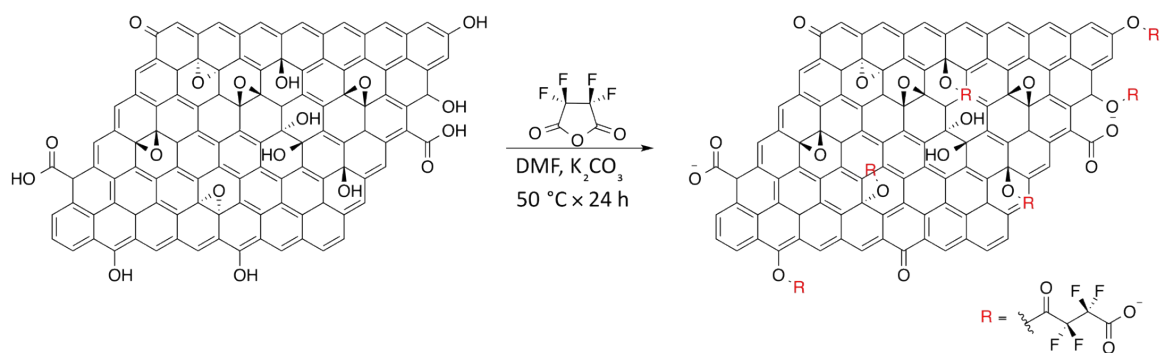
**Table S1.** Position (eV)/FWHM (eV) values of peaks resulting from curve-fitting of C 1s XPS spectra of GO, C-GO, F-GO, C-RGO and RGO.

Sample	C=C	C-OH	C-O-C	C=O + COO <sup>-</sup>	COOH	-CH <sub>2</sub> -	-CF <sub>2</sub> -
<b>GO</b>	284.8/1.6	286.4/1.0	287.0/1.1	288.1/1.3	289.2/1.6	-	-
<b>C-GO</b>	284.8/1.3	286.6/1.1 <sup>a</sup>	287.1/1.3	288.2/1.2	289.3/1.3	285.4/1.5	-
<b>F-GO</b>	284.8/1.4	286.5/1.2 <sup>a</sup>	287.0/1.2	288.3/1.5	289.2/1.3	-	290.0/1.5
<b>C-RGO</b>	284.8/1.3	286.3/1.3 <sup>a</sup>	286.8/1.3	288.3/1.2	289.5/1.2	285.5/1.4	-
<b>RGO</b>	284.8/1.2	286.3/1.4	286.8/1.5	288.0/1.8	289.0/1.6	-	-

<sup>a</sup>In C-GO, F-GO and C-RGO the C-OH signal is also contributed by the estereal C-O(C=O)- atoms of graphene oxide bound to the succinate moiety, which display the same BE as C-OH.



**Figure S3.** XPS spectrum in the F 1s region of F-GO. Predominant component at 689.0 eV is attributed to -CF<sub>2</sub>- groups in succinate moieties appended to GO sheets. Minor component at 686 eV is compatible with ionic F atoms, its origin is uncertain.



**Figure S4.** Reaction scheme for the synthesis of F-GO.