



Cite this: DOI: 10.1039/c9ta90299f

Correction: Flat-shaped carbon–graphene microcomposites as electrodes for high energy supercapacitors

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DOI: 10.1039/c9ta90299f

rsc.li/materials-a

Correction for 'Flat-shaped carbon–graphene microcomposites as electrodes for high energy supercapacitors' by Gelines Moreno-Fernández *et al.*, *J. Mater. Chem. A*, 2019, 7, 14646–14655.

The authors regret a mistake in Table 1 of the published article. The correct version of the Table 1 is shown below.

The authors also wish to amend the following sentences in the original version of the manuscript:

- The sentence "The t-plot method was used to calculate the external surface area (S_{EXT}) in the relative pressure range of 0.07–0.25." included in the physicochemical characterization (page 14648) section should be removed.

- The last sentence of page 14648, where it is written "The increment of S_{EXT} ..." should say "The increment of V_{meso} ..."

- The text on line 7, page 14649 which reads "The isotherm of the activated material also presents a marked micropore contribution but, as evidenced by V_{meso} and S_{EXT} values (Table 1) some narrow mesoporosity is generated during activation." should instead read as follows: "The isotherm of the activated material also presents a marked micropore contribution but, as evidenced by V_{meso} values (Table 1) some narrow mesoporosity is generated during activation".

The authors acknowledge Dr Teresa Centeno, INCAR-CSIC (Spain) for detecting these errors.

Table 1 Total pore volume (V_T), mesopore volume (V_{meso}), micropore volume (V_{DR}), BET specific surface area (S_{BET}), DFT specific surface area (S_{DFT}), total micropore surface area (S_{DR}), narrow micropore volume (V_{DR}), and narrow micropore specific surface area (S_{DR})

	N_2^a					CO_2^b		
	V_T ($cm^3 g^{-1}$)	V_{meso} ($cm^3 g^{-1}$)	V_{DR}^c ($cm^3 g^{-1}$)	S_{BET} ($m^2 g^{-1}$)	S_{DFT}^d ($m^2 g^{-1}$)	S_{DR}^e ($m^2 g^{-1}$)	V_{DR}^c ($cm^3 g^{-1}$)	S_{DR}^e ($m^2 g^{-1}$)
ResFa	0.37	0.13	0.24	598	720	551	0.19	632
ResFaGO	0.50	0.11	0.39	948	1150	880	0.22	840
ResFaGO-A	1.12	0.21	0.91	1961	1991	1373	0.47	1122

^a Data obtained from N_2 adsorption–desorption isotherms at -196 °C. ^b Data obtained from CO_2 adsorption isotherms at 0 °C. ^c Data obtained from Dubinin–Raduskevich. ^d Data obtained from the 2D-NLDFT. ^e Data obtained from Dubinin–Raduskevich.

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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