

## CORRECTION

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

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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 ( $S_{\text{EXT}}$ ) 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  $S_{\text{EXT}}$ ..." should say "The increment of  $V_{\text{meso}}$ ..."

- 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  $V_{\text{meso}}$  and  $S_{\text{EXT}}$  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  $V_{\text{meso}}$  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 ( $V_{\text{T}}$ ), mesopore volume ( $V_{\text{meso}}$ ), micropore volume ( $V_{\text{DR}}$ ), BET specific surface area ( $S_{\text{BET}}$ ), DFT specific surface area ( $S_{\text{DFT}}$ ), total micropore surface area ( $S_{\text{DR}}$ ), narrow micropore volume ( $V_{\text{DR}}$ ), and narrow micropore specific surface area ( $S_{\text{DR}}$ )

	$\text{N}_2^a$						$\text{CO}_2^b$	
	$V_{\text{T}}$ ( $\text{cm}^3 \text{g}^{-1}$ )	$V_{\text{meso}}$ ( $\text{cm}^3 \text{g}^{-1}$ )	$V_{\text{DR}}^c$ ( $\text{cm}^3 \text{g}^{-1}$ )	$S_{\text{BET}}$ ( $\text{m}^2 \text{g}^{-1}$ )	$S_{\text{DFT}}^d$ ( $\text{m}^2 \text{g}^{-1}$ )	$S_{\text{DR}}^e$ ( $\text{m}^2 \text{g}^{-1}$ )	$V_{\text{DR}}^c$ ( $\text{cm}^3 \text{g}^{-1}$ )	$S_{\text{DR}}^e$ ( $\text{m}^2 \text{g}^{-1}$ )
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

<sup>a</sup> Data obtained from  $\text{N}_2$  adsorption–desorption isotherms at  $-196^\circ\text{C}$ . <sup>b</sup> Data obtained from  $\text{CO}_2$  adsorption isotherms at  $0^\circ\text{C}$ . <sup>c</sup> Data obtained from Dubinin–Raduskevich. <sup>d</sup> Data obtained from the 2D-NLDFT. <sup>e</sup> 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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