

## 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 ( $S_{EXT}$ ) 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".

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**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

<sup>a</sup> Data obtained from  $N_2$  adsorption–desorption isotherms at  $-196$  °C. <sup>b</sup> Data obtained from  $CO_2$  adsorption isotherms at  $0$  °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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