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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

[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–Radusckevich. [d] Data obtained from the 2D-NLDFT. [e] Data obtained from Dubinin–Radusckevich.

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

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1486 | *J. Mater. Chem. A*, 2020, **8**, 1486

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