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CORRECTION

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Correction: A supramolecular strategy for tuning the energy level of naphthalenediimide: promoted formation of radical anions with extraordinary stability

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Correction for 'A supramolecular strategy for tuning the energy level of naphthalenediimide: promoted formation of radical anions with extraordinary stability' by Qiao Song *et al.*, *Chem. Sci.*, 2015, **6**, 3342–3346.

The authors regret the incorrect calculation of LUMO and HOMO values in the manuscript. It should be LUMO = $(-4.40 - E_1)$ eV against vacuum.

The LUMO and HOMO values in Fig. 4 should be as following.

The corresponding text should also be corrected as "From the calculated LUMO and HOMO energy shown in Fig. 4, the LUMO energy of NDI/CB[7] is 0.47 eV lower than NDI itself, and therefore the LUMO energy level is as low as −4.54 eV."

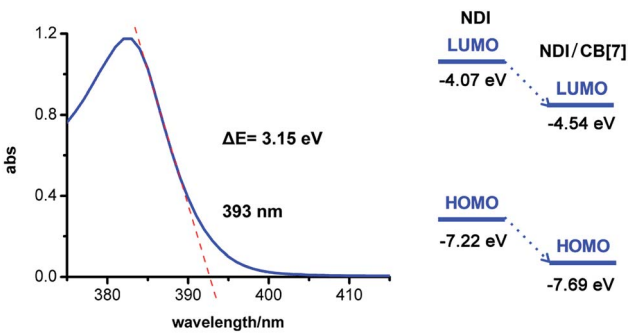

Fig. 4 Calculation of the energy gap between the HOMO and LUMO of NDI and NDI/CB[7] (left); LUMO and HOMO energy of NDI and NDI/CB[7] (right).

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Fig. S10 in ESI should be corrected as:

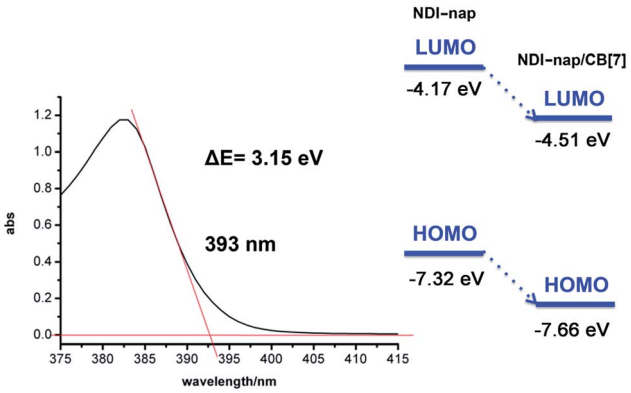

Fig. S10 Calculation of band gap between HOMO and LUMO of NDI-nap and NDI-nap/CB[7] (left); LUMO and HOMO energy of NDI-nap and NDI-nap/CB[7] (right).

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