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CORRECTION

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Correction: Evaluating the role of energetic disorder and thermal activation in exciton transport

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Correction for 'Evaluating the role of energetic disorder and thermal activation in exciton transport' by S. Matthew Menke *et al.*, *J. Mater. Chem. C*, 2016, **4**, 3437–3442.

There is a mistake in the distance dependence of eqn (4), which should read:

$$\Gamma[E, E', d] = \frac{{R_{0,C}}^6}{d^6 \tau_{rad}} \exp\left(-\frac{E_A}{kT}\right) \exp\left(-\frac{E' - E}{kT}\right)$$

There are also errors in Table 1 on page 3440, which affect four callouts from the main text. In Table 1, the second column, "${R_{0,C}}^6/\tau_{rad}$ [nm^6 ns^{-1}]" is actually "${R_{0,C}}/{\tau_{rad}}^{1/6}$ [nm ns$^{-1/6}$]". This correction only changes the second column title and the seventh column values ("$R_{0,C}$") and the corrected table is given below:

Table 1 Model parameters for the KMC simulations predicting the temperature dependence of L_D. Here, ${R_{0,C}}/{\tau_{rad}}^{1/6}$, E_A, and σ are fitting parameters for the KMC simulations. The η_{PL} is separately measured allowing for the tabulation of τ_{rad} and extraction of $R_{0,C}$ from the fit parameters. Shown for comparison is the average molecular separation (d) determined from the thin-film density (ρ) as $d = \sqrt[3]{3/(4\pi\rho)}$

Material	${R_{0,C}}/{\tau_{rad}}^{1/6}$ [nm ns$^{-1/6}$]	E_A [meV]	σ [meV]	η_{PL} [%]	τ_{rad} [ns]	$R_{0,C}$ [nm]	d [nm]
Alq_3	2.3	<10	150	16.2	99	5.0	0.53
DCV3T	1.7	80	33	8.0	17	2.7	0.54
SubPc	1.4	<10	35	1.0	55	2.7	0.48

For consistency, all references to the fitting parameter ${R_{0,C}}^6/\tau_{rad}$ must also be changed to ${R_{0,C}}/{\tau_{rad}}^{1/6}$. This occurs on page 3439, column 1, line 13; page 3439, column 2, line 3; page 3439, column 2, line 18; and page 3440 in the caption of Fig. 4, line 6.

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

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