CORRECTION

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

 Thomas R. Fielitz,[†] S. Matthew Menke[‡] and Russell J. Holmes^{*}

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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.

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

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

There are also errors in Table 1 on page 3440, which affect four callouts from the main text. In Table 1, the second column, " $R_{0,C}^6/\tau_{\text{rad}}$ [$\text{nm}^6 \text{ns}^{-1}$]" is actually " $R_{0,C}/\tau_{\text{rad}}^{1/6}$ [$\text{nm ns}^{-1/6}$]". 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, $R_{0,C}/\tau_{\text{rad}}^{1/6}$, 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	$R_{0,C}/\tau_{\text{rad}}^{1/6}$ [$\text{nm ns}^{-1/6}$]	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 $R_{0,C}^6/\tau_{\text{rad}}$ must also be changed to $R_{0,C}/\tau_{\text{rad}}^{1/6}$. 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.

Department of Chemical Engineering and Materials Science, University of Minnesota, 421 Washington Ave SE, Minneapolis, MN, 55455, USA. E-mail: rholmes@umn.edu

[†] This author identified the errors in the original paper and checked the consistency of the correction with the rest of the paper as published.

[‡] Current address: Department of Physics, Cavendish Laboratory, University of Cambridge, JJ. Thompson Ave, Cambridge, CB3 0HE, UK.