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Author Correction: Efficient electron transfer across hydrogen bond interfaces by proton-coupled and -uncoupled pathways

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Correction to: *Nature Communications* <https://doi.org/10.1038/s41467-019-09392-7>, published online 4 April 2019.

The original version of this Article contained an error in the eighteenth sentence of the third paragraph of the ‘Determination of H_{ab} and k_{ET} data for the Mo_2 dimers’ section of the Results, which incorrectly read ‘For $1a^+$ with $2H_{ab} \gg k_B T$, it is inappropriate to calculate the rate constant using H_{ab} for the PCET reaction in the nonadiabatic regime.’ The correct version states ‘ $2H_{ab} \gg k_B T$ ’ in place of ‘ $2H_{ab} \gg k_B T$ ’.

Also, the third sentence of the Discussion originally incorrectly read ‘The linear relationship of $\ln(k_{ET})$ vs. r_{ab} gives $\beta = 1.25$ (Supplementary Fig. 9).’ The correct version states ‘ R_{ab} ’ instead of ‘ r_{ab} ’.

This has been corrected in both the PDF and HTML versions of the Article.

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