Erratum: Classical Approach to Multichromophoric Resonance Energy Transfer [Phys. Rev. Lett. 115, 110402 (2015)]

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We are grateful to Mr. Simon Axelrod and Dr. Aurelia Chenu for identifying a numerical error in the computer program producing Fig. 2 of original Letter. Correcting this error yielded improved results, as shown here in the figure. Specifically, fitting the data with the functional form $(ae^{-k_f t} + be^{-k_{uf} t} + c)/(a + b + c)$ used in the text yielded constants $a = b = 1$, $c = 2.01$, a fast decay component of $k_{uf} = 13.52 \text{ ps}^{-1}$, and an electronic energy transfer rate of $k_f = 1.11 \text{ ps}^{-1}$. The latter, compared to the value of 0.92 ps^{-1} reported in the text, is seen to be in even better agreement with the experimental value of 1.25 ps^{-1} .

FIG. 2. Normalized total energy within the B800 ring (donor) in LH2 initially excited with a delta pulse. Dashed curve is a best fit to the oscillatory (cyan) result.