Erratum: Tensor Network Simulation of Non-Markovian Dynamics in Organic Polaritons [Phys. Rev. Lett. 121, 227401 (2018)]

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The reorganization energy of the molecular model used in the Letter is actually $\Delta = 35.6$ meV (a factor of π smaller than stated). The necessary corrections are as follows:

(1) The last sentence in the left column on page 2 should read: "For Rhodamine 800 this density is extracted from the spectroscopic measurements in Ref. [34] […] and reorganization energy $\Delta = \int_0^\infty [J_v(\omega)/\pi\omega]\mathrm{d}\omega \approx 35.6$ meV."
(2) In the last paragraph of the right column on page 4, the correct value for τ_{RC} is 23 fs.
(3) The first paragraph in page 5 should read "In addition to the vibration-free polaritons $|\pm\rangle$, emission bands at intermediate energies are visible in the spectrum. This is interpreted as due to small cavity admixtures to dark states, in line with experimental observations [65,72]." We previously misidentified one of the bands as being at the bare-molecule emission frequency, $\omega_e - 2\Delta$, which does not apply with the corrected value of Δ.

We have also updated the Supplemental Material to correct the issues related with the error in the value of Δ. None of the conclusions in the main text or supplemental material are affected.

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