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Addition/Correction

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Correction to Carbon Nanohoops: Excited Singlet and Triplet Behavior of [9]- and [12]-Cycloparaphenylene

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In the abstract and main text of this manuscript, we inadvertently entered wrong estimates of the rate constants of energy transfer (k_{et}) to biphenyl as well as the rate constants of triplet state quenching by molecular oxygen. The correct k_{et} values for reaction 5 are 1.9×10^4 M^{-1} s^{-1} and 1.7×10^4 M^{-1} s^{-1} for [9]CPP and [12]CPP, respectively. The corrected rates of triplet state quenching by oxygen are 1.7×10^9 M^{-1} s^{-1} and 1.8×10^9 M^{-1} s^{-1} for [9]CPP and [12]CPP, respectively. Figure 6 is revised to show these corrected values.

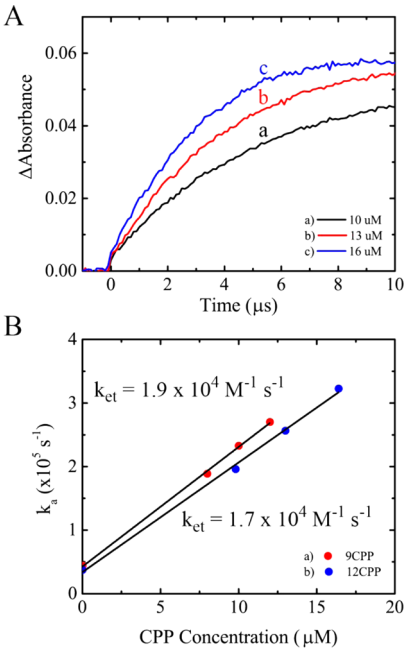

Figure 6. The average lifetime of the nanohoop/oxygen solution with different concentrations of dissolved oxygen. The filled dot represents the experimental data of [9]CPP, and the hollow dot represents [12]CPP. The experimental trends are fit to a straight line (red = [9]CPP; blue = [12]CPP), and the slope of the line is taken to be the bimolecular reaction rate of triplet quenching by oxygen.

Table 2 is revised to reflect the corrected values, and eq 1 is revised to reflect correct subscripts.

$$\Phi_f = \Phi_{std}(F_{nh}/F_{std})(1 - 10^{Astd}/1 - 10^{Anh})(n^2{}_{nh}/n^2{}_{std}) \quad (1)$$

Table 2. Summary of Measured Extinction Coefficients and Φ Values

sample	λ_{max} (nm)	ε_T (M^{-1} cm^{-1})	k_{et} (M^{-1} s^{-1})	Φ_f	Φ_T	Φ_{nr}
[9]CPP	390	25000 ± 4000	1.9×10^4	0.46	0.18	0.36
[12]CPP	680	31000 ± 1300	1.7×10^4	0.83	0.13	0.04

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