Erratum: Binding energies of trions and biexcitons in two-dimensional semiconductors from diffusion quantum Monte Carlo calculations [Phys. Rev. B **95, 081301(R) (2017)]**

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We have noticed a typographical error in the definition of the r_* parameter below Eq. (2). The definition should be $r_* = 2\pi\kappa_{\perp}/\epsilon$. This change does not affect any of our results. To clarify, the values reported in Table I were calculated for free-standing monolayers for which $\epsilon = 1$.