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Erratum: "Semilocal and hybrid density embedding calculations of ground-state charge-transfer complexes" [J. Chem. Phys. 138, 124112 (2013)] **FREE**

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Erratum: “Semilocal and hybrid density embedding calculations of ground-state charge-transfer complexes” [J. Chem. Phys. **138**, 124112 (2013)]

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In Table I of Ref. **1** the error on embedding energy (ΔE) is reported with inverted sign for all the semilocal and hybrid exchange-correlation approximations.

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