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Erratum

E2 and S_N2 Reactions of X^- + CH_3CH_2X (X = F, Cl); an *ab Initio* and DFT Benchmark Study. [*J. Chem. Theory Comput. 4,* 929–940 (2008)]. By A. Patrícia Bento, Miquel Solà, and F. Matthias Bickelhaupt*.

Page 933. In Table 1, the CCSD(T) value of the **1bPC** species at the CBS limit obtained from two-point fits (aug-cc-pVTZ and aug-cc-pVQZ) is −34.27 kcal/mol and not −37.39 kcal/mol as indicated in the original paper. Conclusions are not affected, because they are based on the aug-cc-pVQZ values and not on the CBS extrapolations.

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