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Erratum

Erratum to "He I photoelectron spectroscopy of formic acid isotopomers HCOOH and DCOOD" [Chem. Phys. 272 (2001) 77–90]☆

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Due to a computer word replacement slip up in the submitted manuscript, the word "*cis* (or *anti*)" and "*trans* (or *syn*)" should be interchanged in the first paragraph of page 79. This paragraph should now read "The ground state of neutral formic acid is planar and belongs to the C_S symmetry group, according to the geometrical structure determined by microwave spectroscopy of a large set of formic acid isotopomers. Its structure depicted in Fig. 3, from the work of Davis et al. [9], corresponds to the lowest energy isomer, of *trans* (or *syn*) configuration. The *cis* (or *anti*) isomer of HCOOH lies about 90 meV above the *trans* isomer according to gas phase infrared spectroscopic studies [10] and more precisely at 169 meV, from the assignment and relative intensities of weak lines in the microwave spectrum [11,12]".

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