

Erratum: Theoretical characterization of the potential energy surface for $\text{H}+\text{O}_2 = \text{HO}_2^* = \text{OH}+\text{O}$. III. Computed points to define a global potential energy surface [J. Chem. Phys. **94, 7068 (1991)]**

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At several points in Sec. IV of the paper, and in the titles and footnotes of the Tables, the basis set utilized for most of the CASSCF/CCI calculations described in this paper was incorrectly identified as $[4s3p2d1f/3s*2p]$. The correct designation is $[4s3p2d1f/3s*2p1d]$.

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