ADDITIONS AND CORRECTIONS

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W. Q. Tian, M. Ge,* B. R. Sahu, and D. Wang: Geometrical and Electronic Structure of Pt_7 Cluster: A Density Functional Study

Figure 4c in the original article contained an error. The correct figure is Figure 1 here.

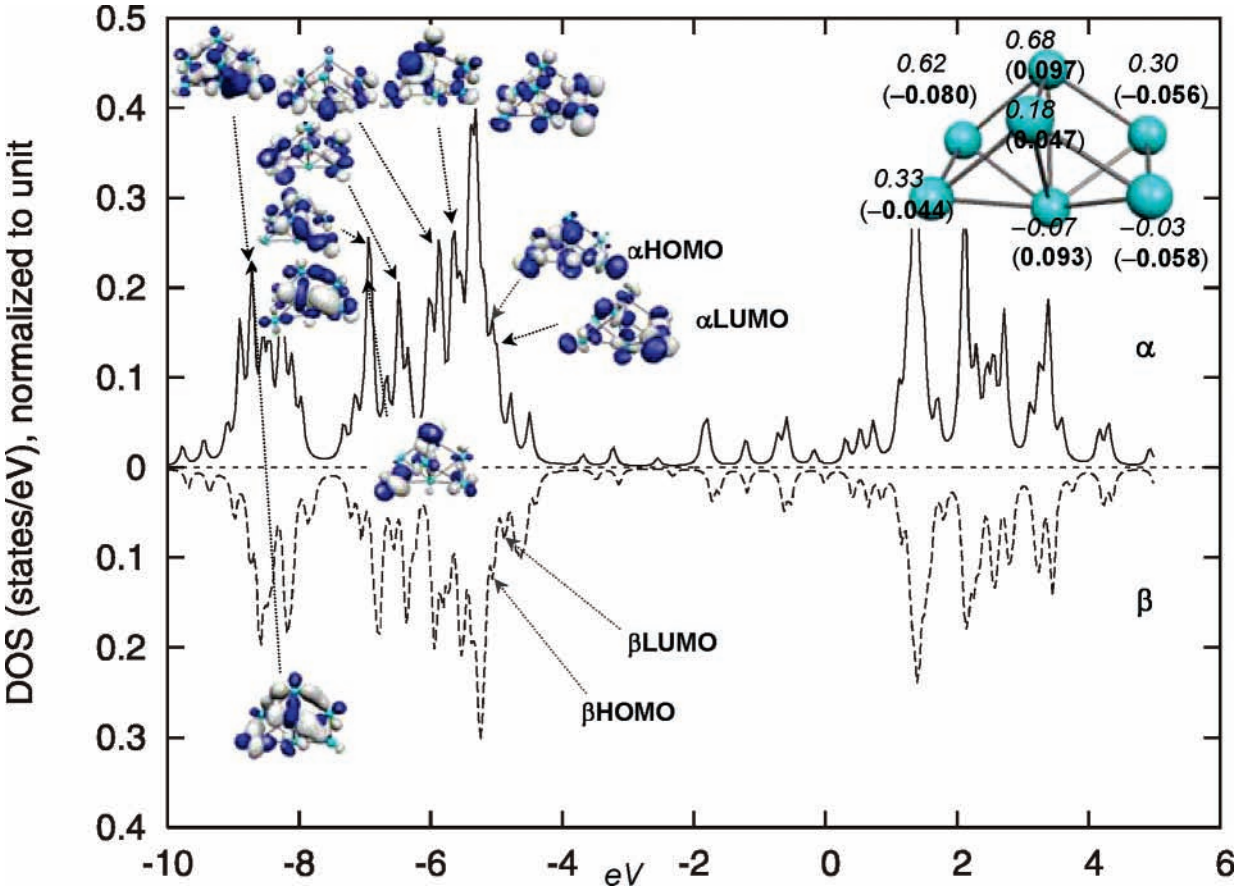

Figure 1. Figure 1. (replacement for Figure 4c in the original article). The α- and *β*- spin density of states of triplet CTP Pt_7 cluster. The net spin charges and the natural atomic orbital partial charges (in parentheses) are also shown.

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