Editor's Comment:

The authors give no indication with regard to the motivation of their work. There are many computer calculations of the electronic structure of carbon allotropes. Why did these authors undertake a new one? Do their results differ from others' results, do their results agree? The ground state energies should be given per molecule, 0% error is
not professional.

Until the authors do not explain their motivation, the status of their results and do not correct the errors, I do not recommend publication.

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