The XXI International Scientific Conference of Young Scientists and Specialists (AYSS-2017)



Contribution ID: 259 Type: Oral

Neon three - atomic clusters within Faddeev approach

This work is aimed at a theoretical investigation of the neon atomic clusters. Bound state energies and wave functions have been calculated. The modern realistic potential models are used for description the interatomic interaction. Differential Faddeev equations in the total angular momentum representation are used for the binding energies calculations of neon trimer system. The results obtained are compared with previous published results.

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Track Classification: Theoretical Physics