

Electron Correlation Dynamics In Atomic Collisions (Cambridge Monographs On Atomic, Molecular And Chemical Physics) By J. H. McGuire

By J. H. McGuire

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collisions are calculated using time-dependent density functional molecular dynamics and TDDFT are also well for atomic collisions,

"Johnson R E. Introduction to Atomic and Molecular Collisions "McGuire J.H. Electron correlation dynamics in atomic chemistry and chemical physics

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