

Adapting the Ab-Initio IMSRG for Open-Shell Atomic Systems

Wednesday, 10 February 2021 13:45 (15)

The nuclear charge distribution and nuclear magnetic moment modify the Coulomb potential in atoms, resulting in shifts in the electronic levels. Thus, atomic spectroscopy provides a way to probe nuclear structure. These measurements, however, require precise calculations of isotope shift factors and hyperfine constants. The IMSRG is an ab-initio technique, successfully used in nuclei, that evolves a many-body Hamiltonian using continuous unitary transformations. I will present a new application of the IMSRG to atomic systems for calculating spectra and isotope shift factors. I will discuss first results and the current status of these calculations as well as what we hope to achieve moving forward.

email address

tsgaurav@gmail.com

Please select: Experiment or Theory

Theory

Primary author(s) : TENKILA, Gaurav (University of British Columbia)

Presenter(s) : TENKILA, Gaurav (University of British Columbia)