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Frontiers in Ab-Initio Computations of Atomic Nuclei

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Atomic nuclei exhibit multiple energy scales ranging from hundreds of MeV in binding energies to fractions of an MeV for low-lying collective excitations. Describing these different energy scales within an ab-initio framework is a long-standing challenge that we overcome by using high-performance computing, many-body methods with polynomial scaling, and ideas from effective-field-theory. With the recent advancements of ab-initio methods we can now address how collectivity and shape coexistence emerge in nuclei from chiral interactions. We accurately describe the first $2+$ and $4+$ energies and the quadrupole transitions from the first $2+$ to the ground-state in neon isotopes. For $^{32,34}\text{Ne}$ less is known and we predict that they are strongly deformed and collective. For ^{30}Ne we interestingly find that a deformed and nearly spherical shape coexist, similar to what is seen in ^{32}Mg . We also confirm that ^{78}Ni has a low-lying rotational band, and that deformed ground states and shape coexistence emerge along the magic neutron number $N = 50$ towards the key nucleus ^{70}Ca . On the neutron-deficient side we also addressed structure of nuclei around the strongly deformed $N = Z = 40$ nucleus ^{80}Zr , although there are challenges our results are competitive with mean-field calculations. We also made predictions for the magnetic dipole transition in ^{48}Ca . Here we found that the transition strength is consistent with a (γ, n) experiment but is larger than the results from inelastic electron- and proton-scattering experiments. With this talk I hope to convey that the accurate computation of multiscale nuclear physics demonstrates the predictive power of modern ab initio methods.

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Email Address

hageng@ornl.gov

Presenter if not the submitter of this abstract

Primary author: HAGEN, Gaute

Presenter: HAGEN, Gaute

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