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Strong Variation of Density of States and Anomalous Isotope Effect in Low and High Tc Superconductors GUANG-LIN ZHAO, Southern University and A&M College — In this work, first-principles density functional theory (DFT) calculations of electronic structures are integrated into the fundamental formalism of many-body physics for superconductivity in Zr, Nb₃Sn, and YBa₂Cu₃O₇. It is shown that the electronic structures of the transition metals and compounds such as Zr, Nb₃Sn, and YBa₂Cu₃O₇ are very complex. The electron densities of states around their Fermi levels possess sharp variations that have a large contribution to the anomalous isotope effect in these superconductors. The work was funded in part by NSF LASIGMA Project (Award No. EPS-1003897, NSF92010-15-RII-SUBR), AFOSR (FA9550-09-1-0367), and NSF project CBET-0754821.

Guang-Lin Zhao
Southern University and A&M College

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