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In most solid state physics research groups today powerful desktop computers and sophisticated, user-friendly electronic structure codes (ESC) based on density functional theory (DFT) enable students to study complex physical systems with ease. During the course of their study these students learn about DFT and how to use ESC, but seldom get an opportunity to understand the algorithms that make these codes accurate and efficient. As a result they do not develop the necessary computational skills that make them transferable to other scientific working environments. In an attempt to address this, we have developed paradigms to the electronic structure problem that involve self-contained mini computational problems that enhance students' understanding of particular key aspects of ESC. In this talk we give an overview of the mini projects developed thus far as well as those we will work on in future.

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