- C5) Alkali metal atoms have one electron outside a closed shell. When these form solids, assume that the kinetic energy of these valence electrons can be calculated in a free electron approximation, and that the (Coulomb) potential energy of the *whole system* is approximately inversely proportional to the nearest neighbor spacing between the nuclei which form a body centered cubic structure. Define what system of units you are using; in choosing, you may wish to consider that you will be asked for final answers in Rydberg energy units, and Bohr length units.
- a) (2 points) Find the kinetic energy per unit volume in terms of the Fermi wave vector $k_{\rm F}$.
- b) (2 points) Find the number of valence electrons per unit volume in terms of $k_{\rm F}$, and thereby find the kinetic energy per electron.
- c) (2 points) Using the assumption in the preamble above, write the potential energy as a function of $k_{\rm F}$ and an unknown constant α .
- d) (2 points) Suppose we use the experimentally known density of the equilibriumbulk solid (at zero absolute temperature) and determine that $k_{\rm F}=0.5~{\rm Bohr^{-1}}$. Use this information to predict the total energy per electron in Rydberg energy units.
- e) (2 points) Suppose it is known experimentally that the ionization energy of a single atom of this material is 0.3 Rydberg. Use this fact to discuss the accuracy of your answer to part d).