Exercise Sheet 6

1. Exchange-correlation potential

Read the article Parameterised local spin density exchange-correlation energies and potentials for electronic structure calculations I: Zero temperature formalism by J.M. MacLaren, D.P. Clougherty, M.E. McHenry, and M.M. Donovan in Computer Physics Communications **66**, 383 (1991) and implement the Vosko-Wilk-Nusair parametrization for the correlation potential assuming $n_{\uparrow} = n/2 = n_{\downarrow}$ (paramagnetic electron density).

2. LDA calculations with spherical potentials

In the spherical-potential approximation we assume that the charge density coming from orbitals with l > 0 are spherically symmetric (as those with l = 0 are). Using this assumption, perform a self-consistent LDA calculation for He in $1s^2$ configuration.