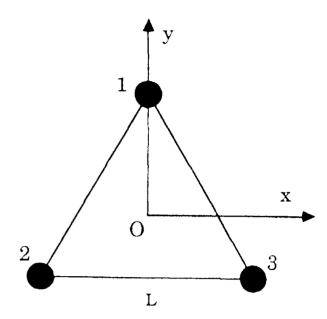
A molecule consists of three equal atoms which form an equilateral triangle of side L as shown. The eigenstates of an electron in this molecule can be considered as linear combinations of orthonormal basis states,  $\phi_i$ , centered on each atom i=1,2,3. In this basis the Hamiltonian has diagonal matrix elements  $\langle \phi_i | H | \phi_i \rangle = \varepsilon$ , and off-diagonal matrix elements between neighboring basis states i and j,  $\langle \phi_i | II | \phi_j \rangle = -t$ , where the energy t > 0.



(a) Define the operator R which operates on the basis functions as follows:

 $R\phi_i = \phi_{i+1}, i = 1, 2; R\phi_3 = \phi_1$ . Note that R has the effect of rotating the basis functions around the triangle and that  $R^3 = 1$ . Show that R commutes with the Hamiltonian.

(b) The eigenstates  $\psi_n$  of R with eigenvalues  $\lambda_n$  can be expressed as  $\psi_n = \sum_{i=1}^3 a_{ni} \phi_i$ . Find the expansion coefficients  $a_{ni}$  and the eigenvalues  $\lambda_n$ .

[Hint: note that  $\lambda_n^3 = 1$ .]

- (c) Give the energy eigenvalues in terms of  $\varepsilon$  and t and give their degeneracies.
- (d) Now suppose an electric field of strength F is switched on in the  $\pm x$ -direction.

The only effect is to change the diagonal matrix elements of the Hamiltonian at each site by an amount  $eFx_i$ , where  $x_i$  is the x position of the center of atom i, and -e is the charge of the electron. Find the new energy eigenvalues to linear order in F.