

Consider a system of electrons with the following eigenstates.

$$\begin{aligned} |0\rangle &= (1, 0, 0, 0) && \text{no electrons} \\ |1\rangle &= (0, 1, 0, 0) && \text{one electron in state } \phi_1 \\ |2\rangle &= (0, 0, 1, 0) && \text{one electron in state } \phi_2 \\ |3\rangle &= (0, 0, 0, 1) && \text{two electrons, one in state } \phi_1, \text{ one in state } \phi_2 \end{aligned}$$

Let electron states  $\phi_n$  be modeled by a one dimensional box of length  $L$ .

$$\phi_n(x) = \sqrt{\frac{2}{L}} \sin\left(\frac{n\pi x}{L}\right)$$

Let  $|\xi\rangle$  be a state vector.

$$|\xi\rangle = c_0|0\rangle + c_1|1\rangle + c_2|2\rangle + c_3|3\rangle, \quad \langle\xi|\xi\rangle = 1$$

Let us determine matrix  $\hat{E}$  such that the expected energy  $\langle E \rangle$  is

$$\langle E \rangle = \langle \xi | \hat{E} | \xi \rangle$$

Matrix  $\hat{E}$  is the sum of kinetic energy  $\hat{K}$  and potential energy  $\hat{V}$ .

$$\hat{E} = \hat{K} + \hat{V}$$

Matrix  $\hat{K}$  can be determined from energy eigenvalues of the box model.

$$\hat{K} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & E_1 & 0 & 0 \\ 0 & 0 & E_2 & 0 \\ 0 & 0 & 0 & E_1 + E_2 \end{pmatrix}, \quad E_n = \frac{n^2 \pi^2 \hbar^2}{2mL^2}$$

Matrix  $\hat{V}$  has one entry due to Coulomb interaction in the two electron state.

$$\hat{V} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & V \end{pmatrix}$$

Let  $\psi(x, y)$  be the antisymmetrized wavefunction of the two electrons.

$$\psi(x, y) = \frac{1}{\sqrt{2}} (\phi_1(x)\phi_2(y) - \phi_1(y)\phi_2(x))$$

Then

$$V = \frac{e^2}{4\pi\epsilon_0} \int_0^L \int_0^L \psi^*(x, y) \left( \frac{1}{|x - y|} \right) \psi(x, y) dx dy$$

Let us now choose  $L = 10^{-9}$  meters and compute numerical values.

$$\hat{K} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.38 \text{ eV} & 0 & 0 \\ 0 & 0 & 1.50 \text{ eV} & 0 \\ 0 & 0 & 0 & 1.88 \text{ eV} \end{pmatrix}$$

Computing  $V$  by numerical integration we have

$$\hat{V} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & 4.67 \text{ eV} \end{pmatrix}$$

Hence

$$\hat{E} = \hat{K} + \hat{V} = \begin{pmatrix} 0 & 0 & 0 & 0 \\ 0 & 0.38 \text{ eV} & 0 & 0 \\ 0 & 0 & 1.50 \text{ eV} & 0 \\ 0 & 0 & 0 & 6.55 \text{ eV} \end{pmatrix}$$

The expected energy is

$$\langle E \rangle = \langle \xi | \hat{E} | \xi \rangle = 0.38 |c_1| + 1.50 |c_2| + 6.55 |c_3|$$

For a state  $|\xi\rangle$  with uniform probability distribution  $|c_i| = \frac{1}{4}$  we have

$$\langle E \rangle = \frac{1}{4}(0.38 + 1.50 + 6.55) = 2.11 \text{ eV}$$