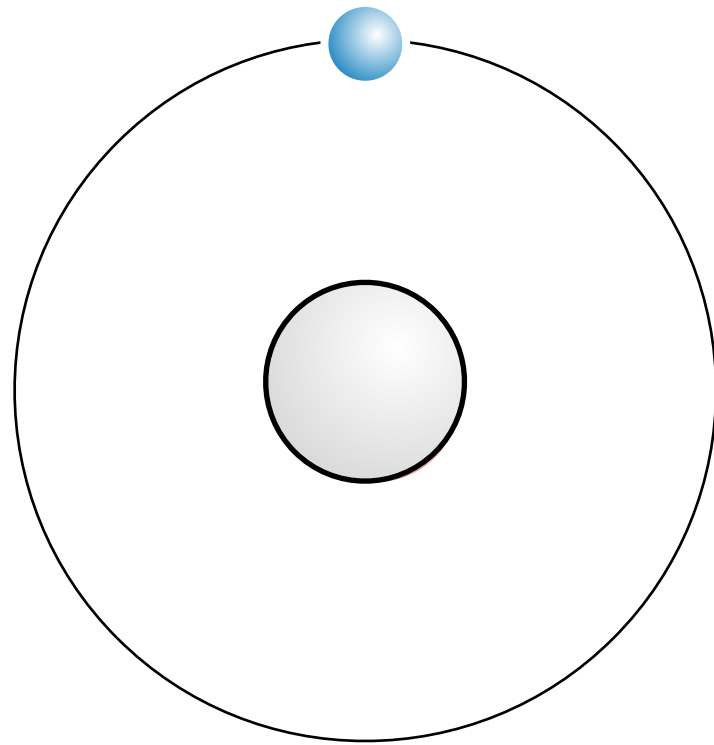
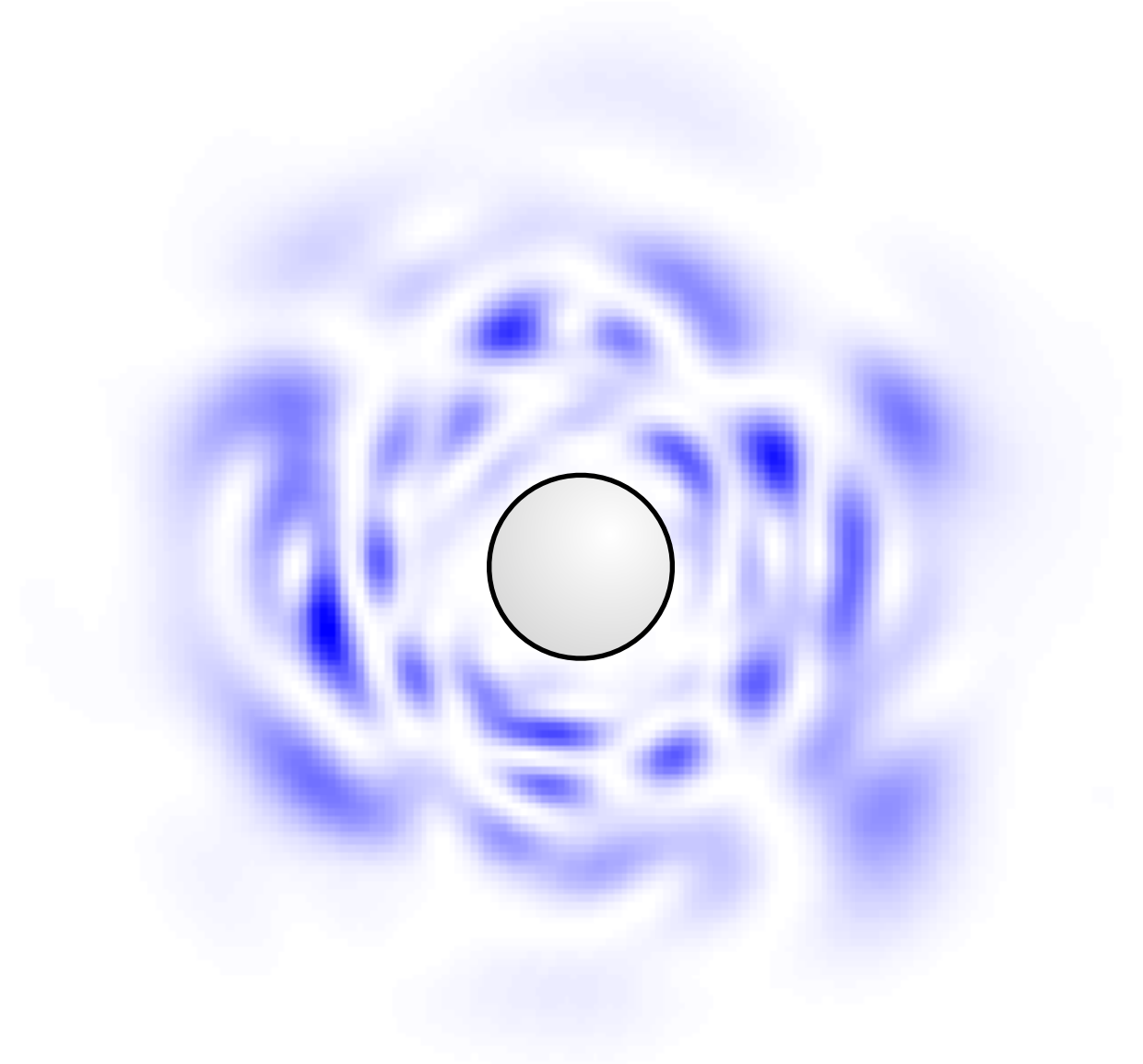


# **Introduction: Schrödinger's Equations**



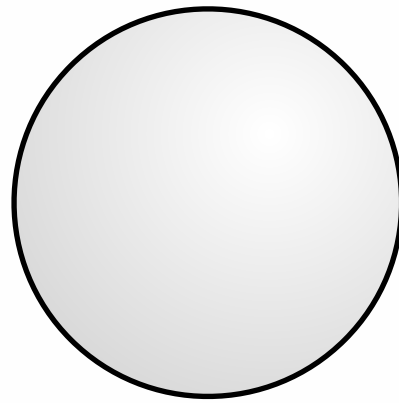
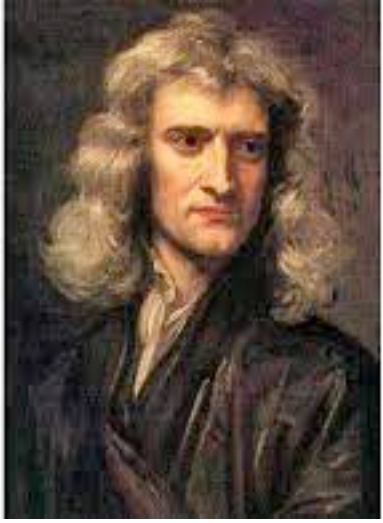
**vs.**



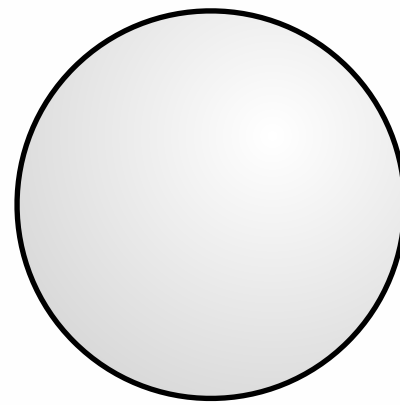
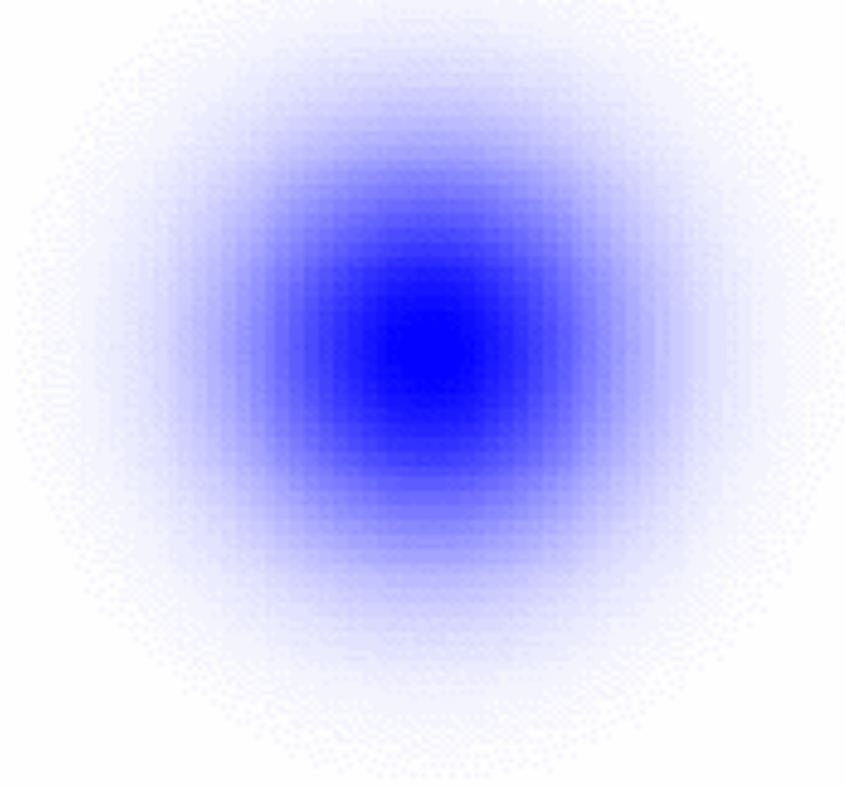
Isaac Newton  
(1643-1727)



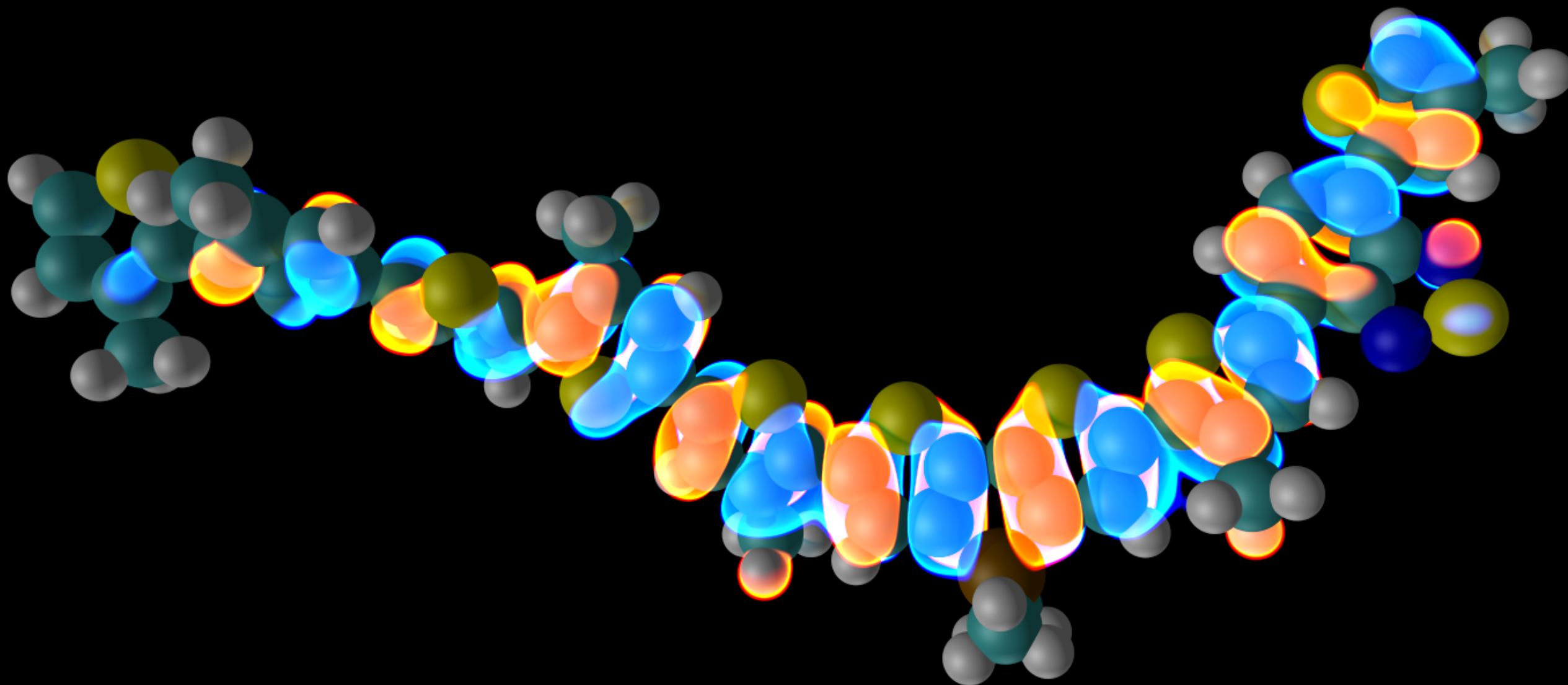
Erwin Schrödinger  
(1887-1961)



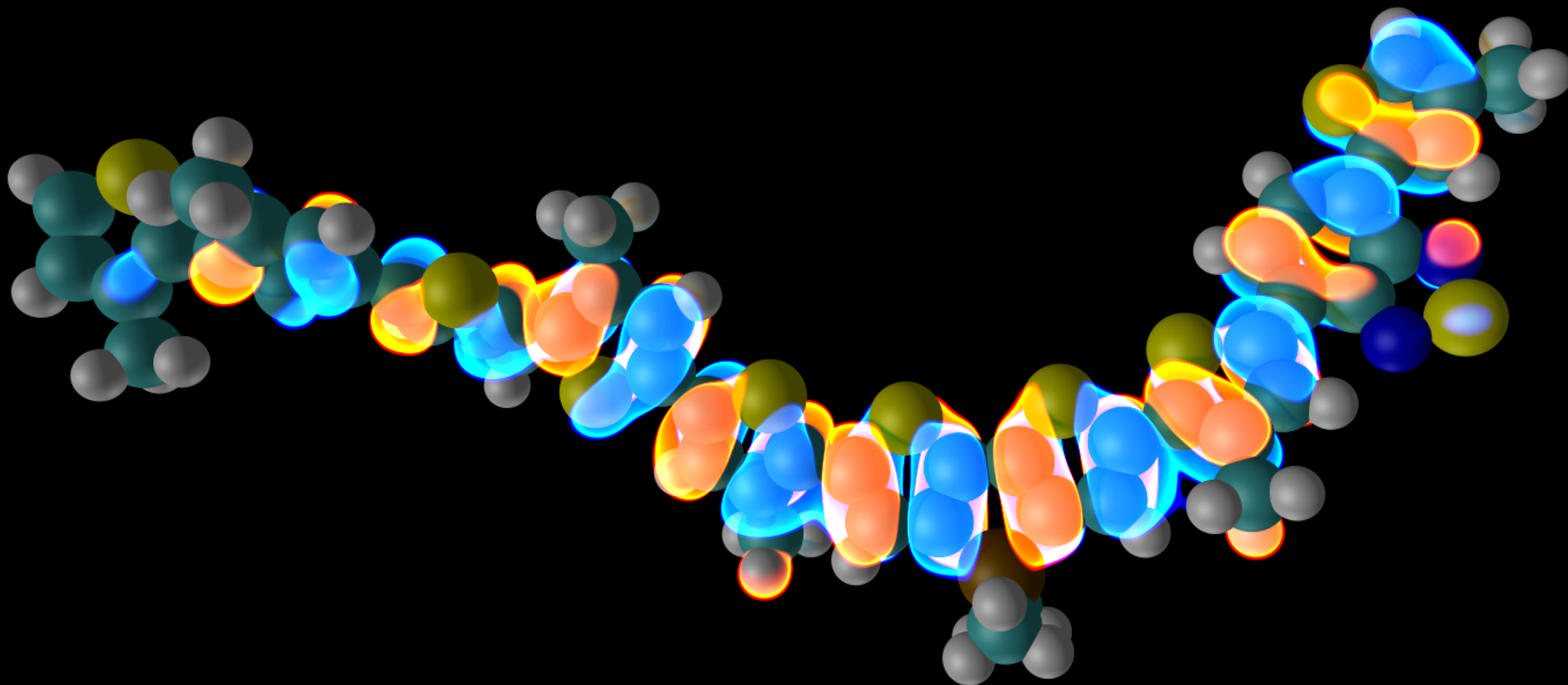
hydrogen atom



hydrogen atom



**material = nuclei + electrons**



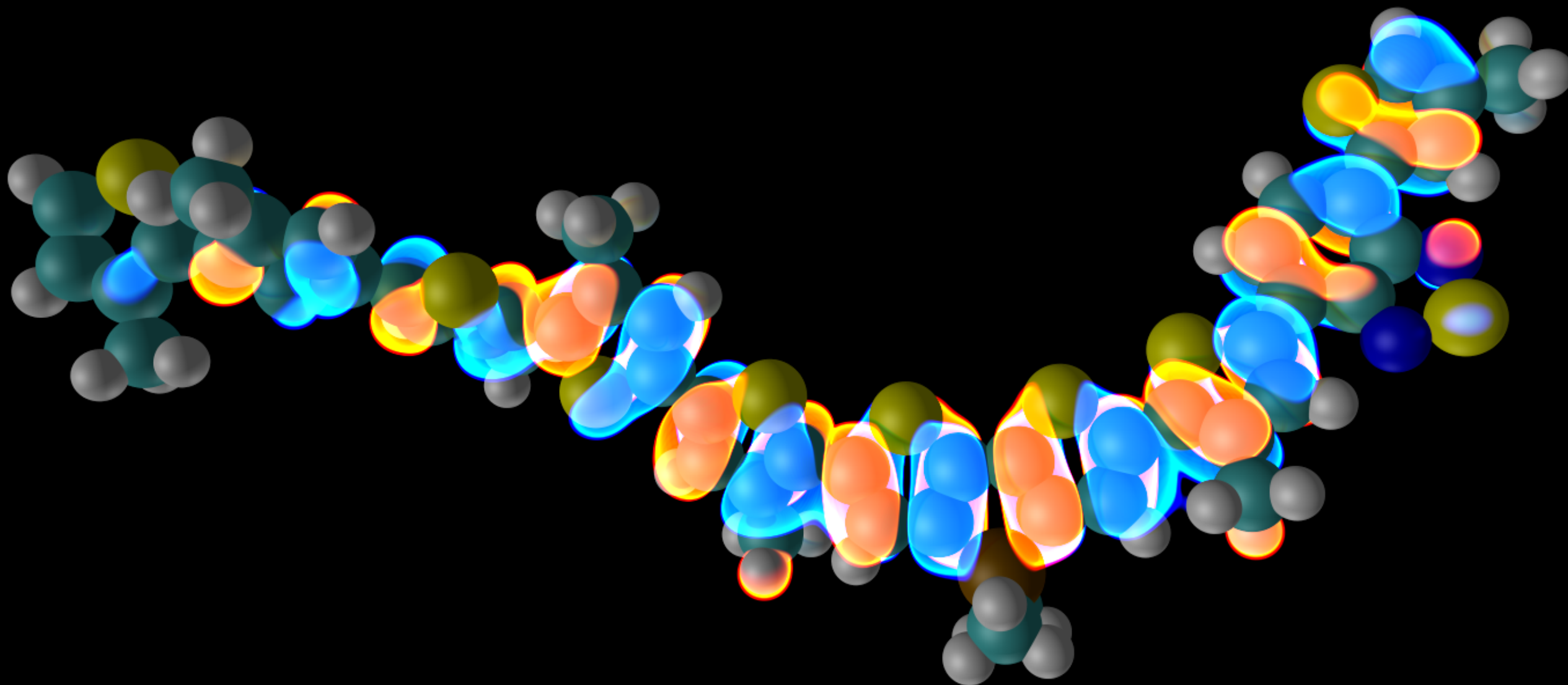
**material = nuclei + electrons**

⎵

**+**

⎵

**-**



**material = nuclei + electrons**

**'glue'**

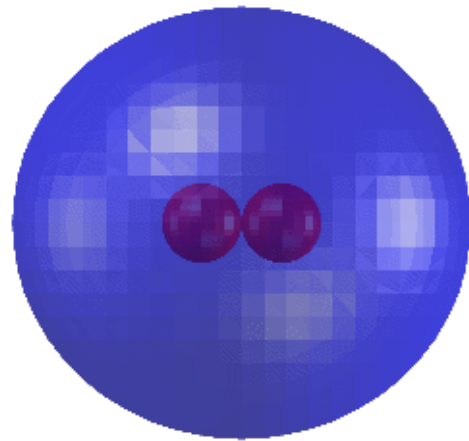




**Schrödinger's equation**

$$H\Psi = E\Psi$$





# Fundamentals

# Classical vs. quantum

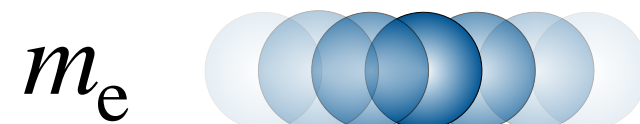
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- **Classical mechanics** and **quantum mechanics** differ in how they describe **electrons**

In classical mechanics, **electrons** are points



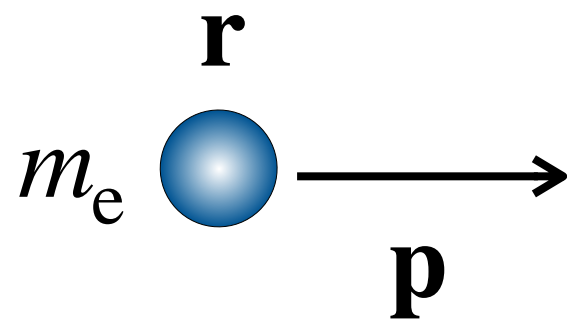
In quantum mechanics, **electrons** are clouds



# Classical vs. quantum

- **Classical mechanics** and **quantum mechanics** differ in how they describe **electrons**

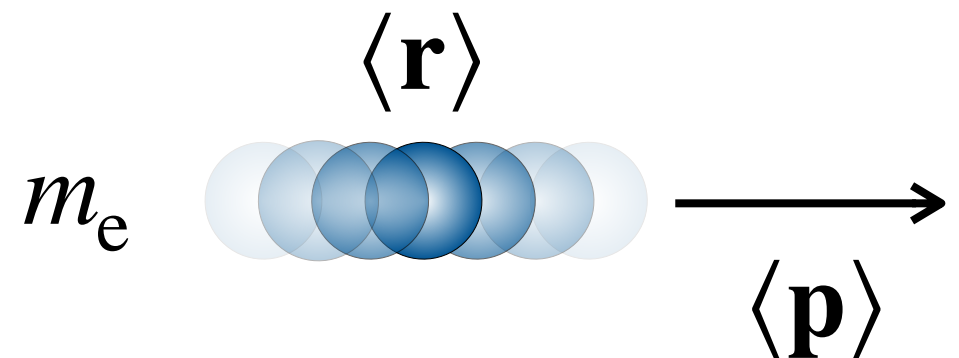
In classical mechanics, the position and momentum are precisely known



In classical mechanics, the position and momentum tell everything about the **electron**

$$(\mathbf{r}, \mathbf{p})$$

In quantum mechanics, only their averages are known



In quantum mechanics, what defines the state of the **electron** is the **wave function**

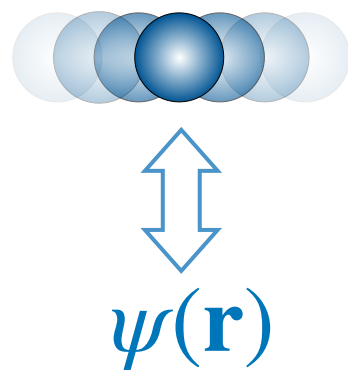
$$\psi(\mathbf{r})$$

# Wave function

- In quantum mechanics, an **electron** is described by a **wave function**, which encodes all information about its state

## Definition | Wave function

The wave function  $\psi(\mathbf{r})$  of an electron is a function of the space variable  $\mathbf{r}$  describing the state of the electron. Its values are complex. Its square modulus gives the probability density, i.e.,  $\rho(\mathbf{r}) = |\psi(\mathbf{r})|^2$ .



A **wave function**  $\psi(\mathbf{r})$  is a function of space whose values are complex and which defines the state of an electron

Probability density

$$\rho(\mathbf{r}) = |\psi(\mathbf{r})|^2 = [\text{Re}(\psi(\mathbf{r}))]^2 + [\text{Im}(\psi(\mathbf{r}))]^2$$

$$\int |\psi(\mathbf{r})|^2 d\mathbf{r} = 1$$

Normalization condition

# Average energy of a wave function

- In quantum mechanics, only the average of the total energy  $\langle E \rangle$  of the wave function is known
- This average energy is also called the expectation value of the total energy

## Definition | Expectation value of the total energy

The **expectation value of the total energy** of an electron is the average value of the total energy over experimental trials for a given state of that electron. If  $\psi(\mathbf{r})$  is the wave function of the electron, this value is

$$\langle E \rangle = \int \underbrace{\psi^*(\mathbf{r})}_{\text{Complex conjugate}} \left( -\frac{\hbar^2}{2m_e} \underbrace{\nabla^2}_{\partial^2/\partial x^2 + \partial^2/\partial y^2 + \partial^2/\partial z^2} \psi(\mathbf{r}) + v(\mathbf{r})\psi(\mathbf{r}) \right) d\mathbf{r}$$

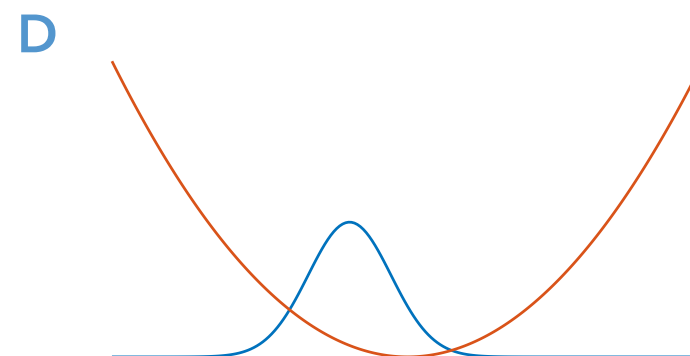
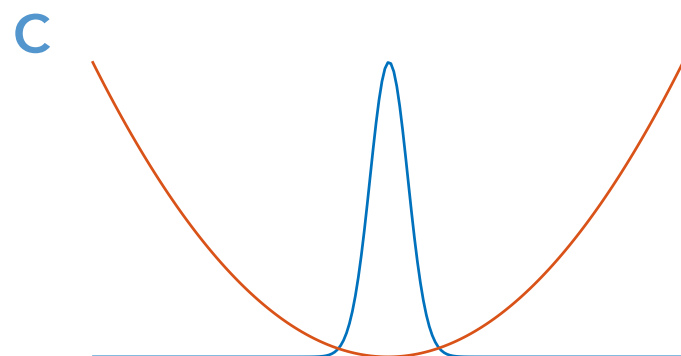
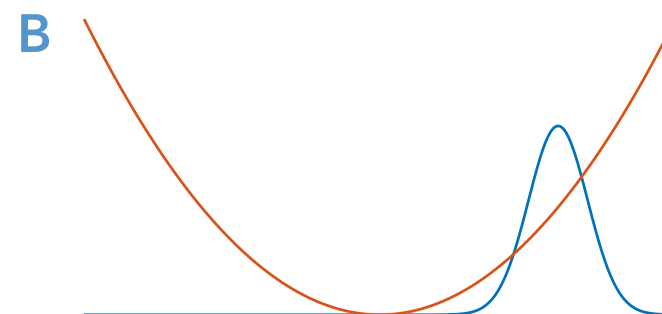
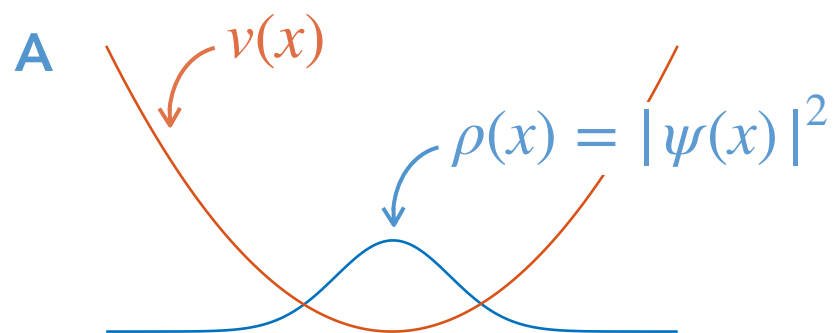
# Understanding each term of the total energy

- The total energy can be split into **kinetic** and **potential** parts

$$\langle K \rangle = \int \psi^*(\mathbf{r}) \left( -\frac{\hbar^2}{2m_e} \nabla^2 \psi(\mathbf{r}) \right) d\mathbf{r} = -\frac{\hbar^2}{2m_e} \int \psi^*(\mathbf{r}) \nabla^2 \psi(\mathbf{r}) d\mathbf{r}$$

$$\langle U \rangle = \int \psi^*(\mathbf{r}) v(\mathbf{r}) \psi(\mathbf{r}) d\mathbf{r} = \int v(\mathbf{r}) \psi^*(\mathbf{r}) \psi(\mathbf{r}) d\mathbf{r} = \int v(\mathbf{r}) \rho(\mathbf{r}) d\mathbf{r}$$

- Problem.** Find the order of  $\langle K \rangle$  and  $\langle U \rangle$  for the 4 cases below:



Increasingly curved  $\psi(x)$

↑

C  
B  
D  
A

Increasingly positive  $v(x)$

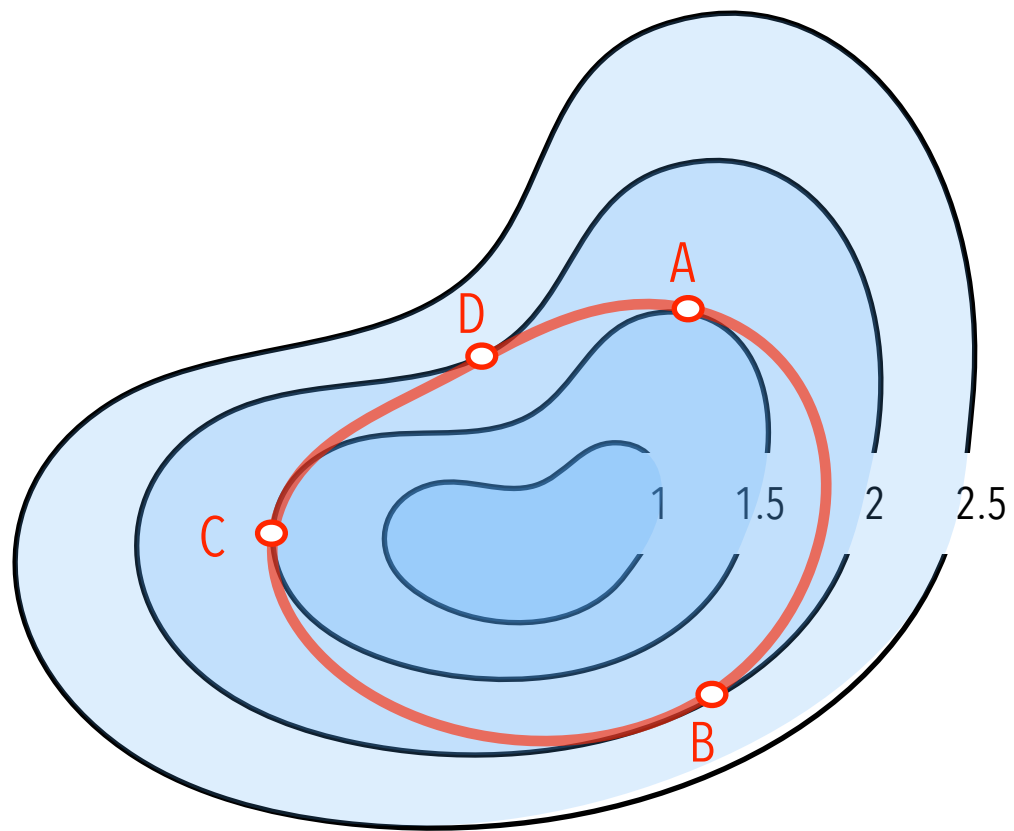
↑

B  
D  
A  
C



# Finding the minimum of the total energy

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# Equilibrium states

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- Using the Lagrange multiplier method, we get the following important result

## Definition | time-independent Schrödinger equation

The equilibrium states  $\psi$  of an electron for a system described by the Hamiltonian  $H$  can be determined by solving the eigenvalue problem

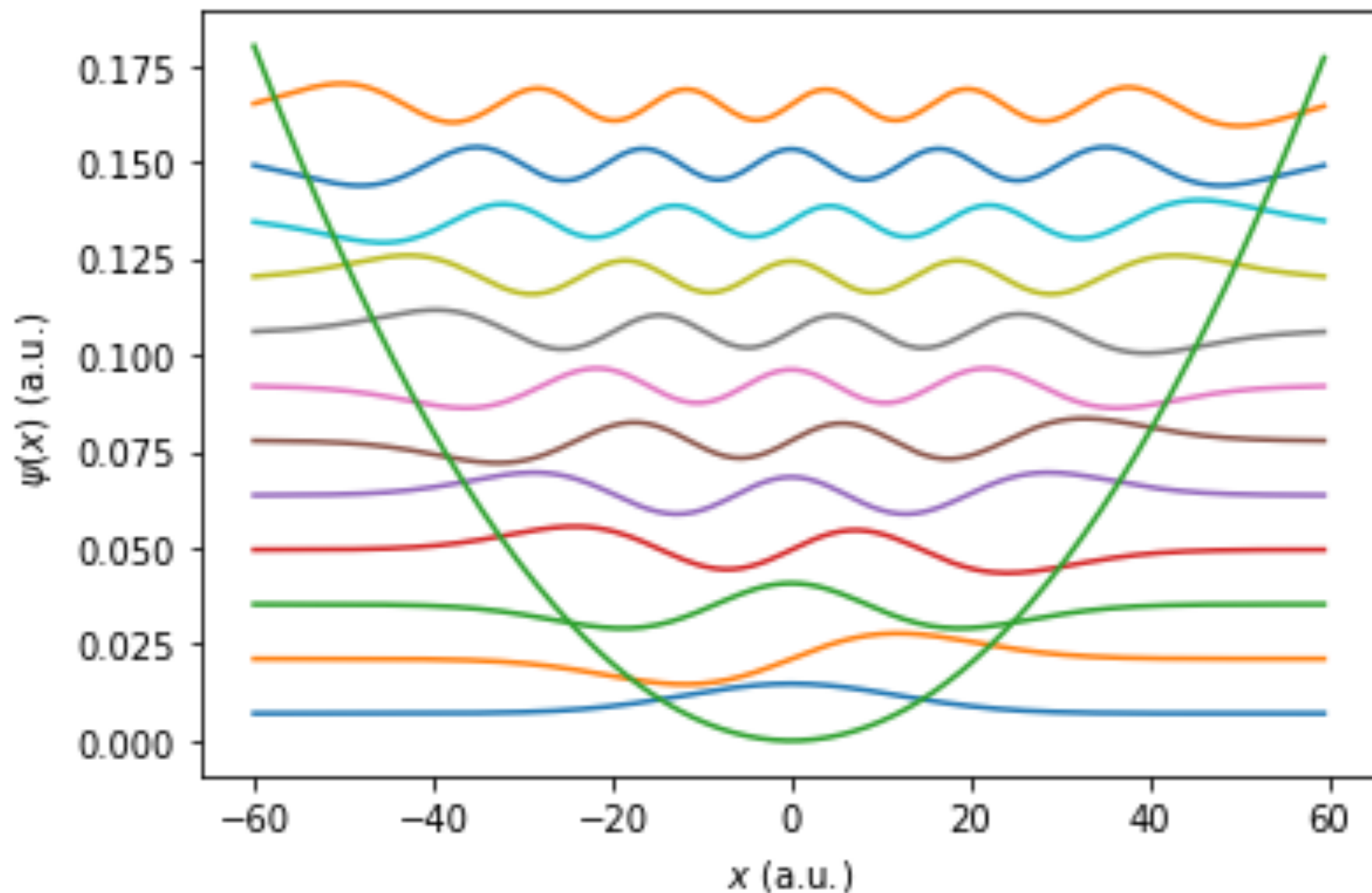
$$H\psi = \mathcal{E}\psi$$

This eigenvalue problem is referred to as the time-independent Schrödinger equation

- **Simulation.** In MATLAB, eigenvalue problems can be solved with `[V,D]=eigs(H,N,'sm')`. Using `electronic_states`, plot the equilibrium states of an electron in a well.

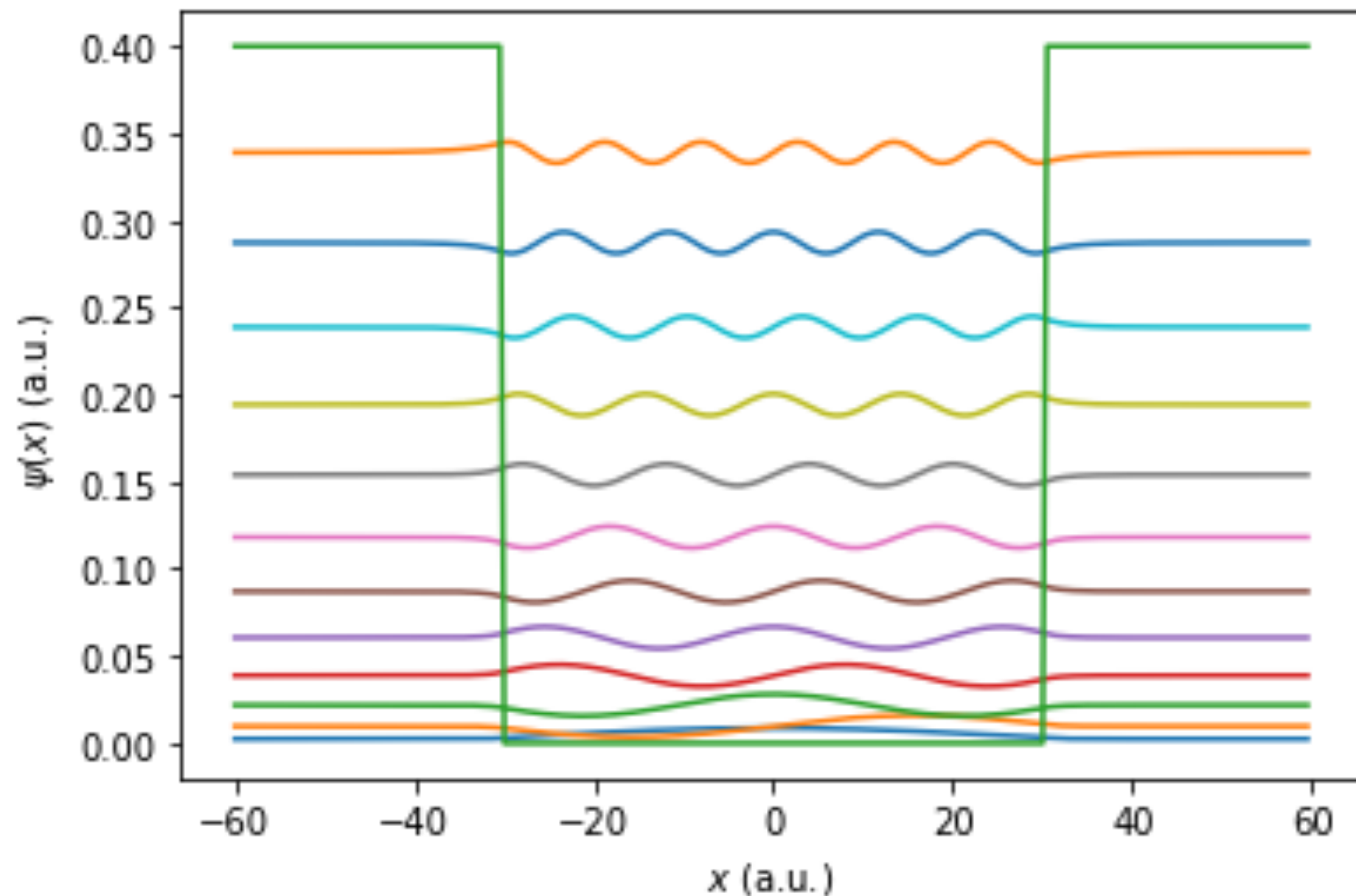
# Equilibrium states (cont'd)

- **Answer.** We obtain a series of quantum states, whose spacing depends on the shape of the well



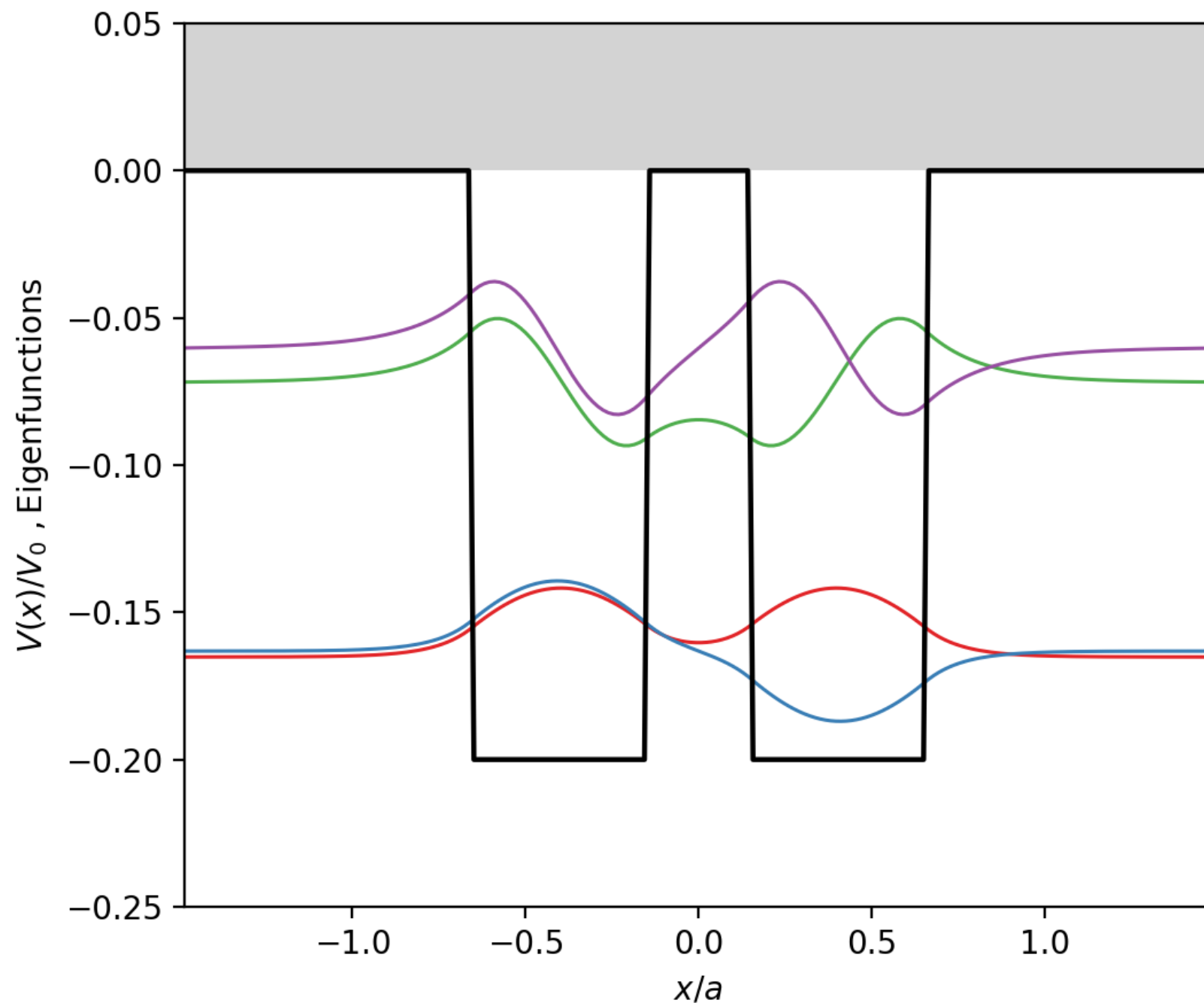
# Equilibrium states (cont'd)

- **Answer.** We obtain a series of quantum states, whose spacing depends on the shape of the well



# Equilibrium states (cont'd)

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