

1. The action principle

- For small $\delta s \in \mathbb{R}$, $f(s + \delta s) = f(s) + \frac{df(s)}{ds} \delta s + R(s, \delta s)$
- With $\delta f := f(s + \delta s) - f(s)$, $\delta f = \frac{df(s)}{ds} \delta s + R(s, \delta s)$, with

$$\lim_{\delta s \rightarrow 0} \frac{R(s, \delta s)}{\delta s} = 0$$

So δf vanishes to first order in δs , so $R(s, \delta s)$ can be written as $O((\delta s)^2)$

- At the extrema of f , $\frac{df(s)}{ds} = 0$ so $\delta f = O((\delta s)^2)$
- **Functional:** map from functions to \mathbb{R}
- $y(t)$ **stationary** for functional S if

$$\frac{dS[y(t) + \varepsilon z(t)]}{d\varepsilon} \Big|_{\varepsilon=0} = 0$$

for every smooth $z(t)$ with $z(a) = z(b) = 0$. We use the notation $\delta y(t) = \varepsilon z(t)$. $y(t)$ is called a **path**.

- **Action principle (variational principle):** paths described by particles are stationary paths of S :

$$\delta S := S[x + \delta x] - S[x] = O((\delta x)^2)$$

for arbitrary smooth small deformations $\delta x(t)$ around true path $x(t)$.

- **Fundamental lemma of the calculus of variations:** Let $f(x)$ be continuous in $[a, b]$ and

$$\int_a^b f(x)g(x) dx = 0$$

for every smooth $g(x)$ in $[a, b]$ with $g(a) = g(b) = 0$. Then $f(x) = 0$ in $[a, b]$.

- **Notation:**

$$\frac{\partial L}{\partial x} = \frac{\partial L(r, s)}{\partial r} \Big|_{(r,s)=(x(t), \dot{x}(t))}, \quad \frac{\partial L}{\partial \dot{x}} = \frac{\partial L(r, s)}{\partial s} \Big|_{(r,s)=(x(t), \dot{x}(t))}$$

- For a path \underline{q} and a Lagrangian $L(\underline{q}, \dot{\underline{q}})$, the action for the path is

$$S = \int_{t_0}^{t_1} L(\underline{q}(t), \dot{\underline{q}}(t)) dt$$

- The action above satisfies

$$0 = \delta S = \int_{t_0}^{t_1} \left(\sum_{i=1}^N \frac{\partial L}{\partial q_i} \delta q_i + \sum_{i=1}^N \frac{\partial L}{\partial \dot{q}_i} \delta \dot{q}_i \right) dt$$

- **Euler-Lagrange equation:**

$$\frac{\partial L}{\partial x} - \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{x}} \right) = 0$$

- The arguments in a Lagrangian, x and \dot{x} , are independent:

$$\frac{\partial x}{\partial \dot{x}} = \frac{\partial \dot{x}}{\partial x} = 0$$

- **Configuration space, \mathcal{C} :** set of all possible instantaneous configurations of a physical system. (Includes positions but not velocities).
- For configuration space \mathcal{C} of system \mathcal{S} , \mathcal{S} has $\dim(\mathcal{C})$ **degrees of freedom**.
- **Generalised coordinates:** A set of coordinates in configuration space.
- Notation: q shows results holds for arbitrary choices of generalised coordinates.
- **Euler-Lagrange equation for configuration space \mathcal{C} :**

$$\frac{\partial L}{\partial q_i} - \frac{d}{dt} \left(\frac{\partial L}{\partial \dot{q}_i} \right) = 0 \quad \forall i \in \{1, \dots, \dim(\mathcal{C})\}$$

- For system with kinetic energy $T(\underline{q}, \underline{\dot{q}})$ and potential energy $V(\underline{q})$, the Lagrangian for the system is

$$L(\underline{q}, \underline{\dot{q}}) = T(\underline{q}, \underline{\dot{q}}) - V(\underline{q})$$

- **Ignorable coordinate q_i :** Lagrangian does not depend on q_i :

$$\frac{\partial L(q_1, \dots, q_N, \dot{q}_1, \dots, \dot{q}_N)}{\partial q_i} = 0$$

- **Generalised momentum** of coordinate q_i :

$$p_i := \frac{\partial L}{\partial \dot{q}_i}$$

- Generalised momentum of ignorable coordinate is conserved.

2. Symmetries, Noether's theorem and conservation laws

- **Transformation depending on ε :** family of smooth maps $\varphi(\varepsilon) : \mathcal{C} \rightarrow \mathcal{C}$ with $\varphi(0)$ the identity map. Can be written as

$$q_i \rightarrow q_i' = \phi_i(q_1, \dots, q_N, \varepsilon)$$

where the ϕ_i are a set of $N = \dim(\mathcal{C})$ functions representing the transformation in the given coordinate system. Change in velocities is

$$\dot{q}_i \rightarrow \frac{d}{dt} \phi_i$$

- **Generator of φ :**

$$\left. \frac{d\varphi(\varepsilon)}{d\varepsilon} \right|_{\varepsilon=0} = \varphi'(0)$$

- In any coordinate system,

$$q_i \rightarrow \phi_i(\underline{q}, \varepsilon) = q_i + \varepsilon a_i(\underline{q}) + O(\varepsilon^2)$$

where

$$a_i = \frac{\partial \phi_i(\underline{q}, \varepsilon)}{\partial \varepsilon} \Big|_{\varepsilon=0}$$

So the generator of the transformation is a_i .

- For velocities,

$$\dot{q}_i \rightarrow \dot{q}_i + \varepsilon \dot{a}_i(q_1, \dots, q_N, \dot{q}_1, \dots, \dot{q}_N) + O(\varepsilon^2)$$

generated by \dot{a}_i .

- Equations of motion don't change when total derivative of function of coordinates and time is added to Lagrangian:

$$L \rightarrow L + \frac{dF(q_1, \dots, q_N, t)}{dt}$$

doesn't change equations of motion.

- Transformation $\varphi(\varepsilon)$ is **symmetry** if for some $F(\underline{q}, t)$,

$$L \rightarrow L' = L(\phi(q_1, \varepsilon), \dots, \phi(q_N, \varepsilon)) = L + \varepsilon \frac{dF(q_1, \dots, q_N, t)}{dt} + O(\varepsilon^2)$$

$F(\underline{q}, t)$ defined up to a constant.

- For ignorable coordinate q_i , transformation $q_i \rightarrow q_i + c_i$ is symmetry since q_i doesn't appear in Lagrangian and \dot{q}_i stays invariant. So $F = 0$ here and $a_k = \delta_{ik}$.
- **Noether's theorem:** Let a symmetric transformation be generated by $a_i(q_1, \dots, q_N)$, so

$$L \rightarrow L + \varepsilon \frac{dF(q_1, \dots, q_N, t)}{dt} + O(\varepsilon^2)$$

Then

$$Q := \left(\sum_{i=1}^N a_i \frac{\partial L}{\partial \dot{q}_i} \right) - F$$

is conserved (so $\frac{dQ}{dt} = 0$).

- Q is called **Noether charge**.
- Given Lagrangian $L(\underline{q}, \underline{\dot{q}}, t)$, **energy** is

$$E := \left(\sum_{i=1}^N \dot{q}_i \frac{\partial L}{\partial \dot{q}_i} \right) - L$$

- Along path $\underline{q}(t)$ satisfying equations of motion,

$$\frac{dE}{dt} = - \frac{\partial L}{\partial t}$$

- So energy conserved iff Lagrangian doesn't depend explicitly on time.

3. Normal modes

- **Canonical** kinetic term: of the form $T = \frac{1}{2} \sum_{i=1}^n \dot{q}_i^2$.

- **Normal mode:** solution to $\ddot{\underline{q}} + A\underline{q} = 0$, associated with eigenvalue $\lambda^{(i)} > 0$ of A , of form

$$\underline{q}(t) = \underline{v}^{(i)} \left(\alpha^{(i)} \cos\left(\sqrt{\lambda^{(i)}} t\right) + \beta^{(i)} \sin\left(\sqrt{\lambda^{(i)}} t\right) \right)$$

- **Zero mode:** solution to $\ddot{\underline{q}} + A\underline{q} = 0$, associated with eigenvalue $\lambda^{(i)} = 0$ of A , of form

$$\underline{q}(t) = \underline{v}^{(i)} \left(\alpha^{(i)} t + \beta^{(i)} \right)$$

- **Instability:** solution to $\ddot{\underline{q}} + A\underline{q} = 0$, associated with eigenvalue $\lambda^{(i)} < 0$ of A , of form

$$\underline{q}(t) = \underline{v}^{(i)} \left(\alpha^{(i)} \cosh\left(\sqrt{-\lambda^{(i)}} t\right) + \beta^{(i)} \sinh\left(\sqrt{-\lambda^{(i)}} t\right) \right)$$

- When no instabilities, general solution is superposition (sum) of normal modes and zero modes.

4. Fields and the wave equation

- **Generalised Euler-Lagrange equations for fields:**

$$\frac{\partial \mathcal{L}}{\partial u} - \frac{\partial}{\partial x} \left(\frac{\partial \mathcal{L}}{\partial u_x} \right) - \frac{\partial}{\partial t} \left(\frac{\partial \mathcal{L}}{\partial u_t} \right) = 0$$

and for n fields $u^{(i)}$:

$$\frac{\partial \mathcal{L}}{\partial u^{(i)}} - \frac{\partial}{\partial x} \left(\frac{\partial \mathcal{L}}{\partial u_x^{(i)}} \right) - \frac{\partial}{\partial t} \left(\frac{\partial \mathcal{L}}{\partial u_t^{(i)}} \right) = 0 \quad \forall i$$

- If fields don't depend on (t, x) but on d coordinates x_i ,

$$\frac{\partial \mathcal{L}}{\partial u^{(i)}} - \sum_{k=1}^d \frac{\partial}{\partial x_k} \left(\frac{\partial \mathcal{L}}{\partial u_k^{(i)}} \right)$$

where $u_k^{(i)} = \frac{\partial u^{(i)}}{\partial x_k}$

- **Massless scalar field Lagrangian:**

$$\mathcal{L} = \frac{1}{2} \rho u_t^2 - \frac{1}{2} \tau u_x^2$$

ρ is **density**, τ is **tension**. The field u is the **massless scalar**.

- Equation of motion for massless scalar field is

$$\rho u_{tt} - \tau u_{xx} = 0$$

which rearranges to **wave equation**:

$$u_{tt} = c^2 u_{xx}$$

where $c^2 = \tau / \rho$.

- **D'Alembert's solution to wave equation:**

$$u(x, t) = f(x - ct) + g(x + ct)$$

$f(x - ct)$ corresponds to a wave moving to the right with speed c , $g(x + ct)$ corresponds to a wave moving to the left with speed c .

- If $u(x, 0) = \varphi(x)$ and $u_t(x, 0) = \psi(x)$ then

$$u(x, t) = \frac{1}{2}(\varphi(x - ct) + \varphi(x + ct)) + \frac{1}{2c} \int_{x-ct}^{x+ct} \psi(s) \, ds$$

- In field theory, **symmetry** is transformation

$$u \rightarrow u' = u + \varepsilon a(u)$$

such that $\delta \mathcal{L} = O(\varepsilon^2)$. $a(u)$ **generates** the transformation.

- **Note:** often, x_0 chosen to be t .
- Let $u_i = \frac{\partial u}{\partial x_i}$, **generalised momentum vector** is

$$\underline{\Pi} := \left(\frac{\partial \mathcal{L}}{\partial u_0}, \dots, \frac{\partial \mathcal{L}}{\partial u_d} \right)$$

- **Noether current** associated to transformation generated by a is

$$\underline{J} = a \underline{\Pi}$$

- If \underline{J} associated to symmetry,

$$\underline{\nabla} \cdot \underline{J} = \sum_{i=0}^d \frac{\partial J_i}{\partial x_i} = 0$$

- **(Noether) charge density:**

$$\mathcal{Q} := J_0$$

- For $d = 1$, **charge contained in interval** (a, b) :

$$Q_{(a,b)} = \int_a^b \mathcal{Q} \, dx$$

- For $d = 1$,

$$\frac{dQ_{(a,b)}}{dt} = J_1(a) - J_1(b)$$

- **Noether charge** is total charge over all space. For $d = 1$:

$$Q := Q_{(-\infty, \infty)} = \int_{-\infty}^{\infty} J_0 \, dx$$

- If $d = 1$ and $\lim_{x \rightarrow \pm\infty} J_1 = 0$,

$$\frac{dQ}{dt} = 0$$

- **Energy-momentum tensor:**

$$T_{ij} := \frac{\partial \mathcal{L}}{\partial u_j} \frac{\partial u}{\partial x_i} - \delta_{ij} \mathcal{L}$$

- **Energy density:**

$$\mathcal{E} := T_{00}$$

- **Conservation law for energy-momentum tensor:**

$$\sum_{j=0}^d \frac{\partial T_{ij}}{\partial x_j} = 0$$

- **Dirichlet boundary condition** for wave equation: $u_t(0, t) = 0$ (so $u(0, t) = 0$ as u has shift symmetry) which gives

$$u(x, t) = f(x - ct) - f(-x - ct)$$

Here, waves reflected off boundary and turned upside down.

- **Neumann (free) boundary condition:** $u_x(0, t) = 0$ which gives

$$u(x, t) = f(x - ct) + f(-x - ct)$$

So waves reflected off boundary and not turned upside down.

- **Junction conditions:**

- u continuous at 0:

$$\lim_{\varepsilon \rightarrow 0^+} u(\varepsilon, t) = \lim_{\varepsilon \rightarrow 0^-} u(\varepsilon, t)$$

- Energy conservation across junction:

$$\frac{d}{dt} \left(\lim_{\varepsilon \rightarrow 0} T(-\varepsilon, \varepsilon) \right) = \lim_{\varepsilon \rightarrow 0} (T_{tx})_{x=-\varepsilon} - \lim_{\varepsilon \rightarrow 0} (T_{tx})_{x=\varepsilon}$$

- **Ansatz for wave function with spring at junction at $x = 0$:**

$$u(x, t) = \begin{cases} \operatorname{Re}((e^{ipx} + R e^{-ipx}) e^{-ipct}) & \text{if } x \leq 0 \\ \operatorname{Re}(T e^{ip(x-ct)}) & \text{if } x > 0 \end{cases}$$

5. The Hamiltonian formalism

- **State** of classical system at given instant in time is complete set of data that fully fixes future evolution of system.
- **Phase (state) space** of system is space of all possible states system can be in at instant in time.
- **Hamiltonian formalism** parameterises phase space as generalised coordinates $\underline{q}(t)$ and associated generalised momenta $\underline{p}(t)$.
- When going from Lagrangian to Hamiltonian formalism, define **generalised momentum** as

$$p_i := \frac{\partial L(\underline{q}, \underline{\dot{q}}, t)}{\partial \dot{q}_i}$$

- **Poisson bracket** of $f(\underline{q}, \underline{p}, t)$ and $g(\underline{q}, \underline{p}, t)$:

$$\{f, g\} := \sum_{i=1}^n \left(\frac{\partial f}{\partial q_i} \frac{\partial g}{\partial p_i} - \frac{\partial f}{\partial p_i} \frac{\partial g}{\partial q_i} \right)$$

where n is dimension of configuration space (half dimension of phase space). Position and momentum treated as independent when taking partial derivatives.

• **Properties of Poisson bracket:**

- **Antisymmetric:** $\{f, g\} = -\{g, f\}$.
- **Linear:** $\{af + bg, h\} = a\{f, h\} + b\{g, h\}$.
- **Leibniz identity:** $\{fg, h\} = f\{g, h\} + g\{f, h\}$.
- **Jacobi identity:** $\{\{f, g\}, h\} + \{\{h, f\}, g\} + \{\{g, h\}, f\} = 0$.
- Let \mathcal{P} be phase space, \mathcal{F} be set of functions from \mathcal{P} to \mathbb{R} .
- **Hamiltonian flow** defined by $f : \mathcal{P} \rightarrow \mathbb{R}$ is infinitesimal transformation on \mathcal{F} given by

$$\Phi_f^{(e)} : \mathcal{F} \rightarrow \mathcal{F}, \quad \Phi_f^{(e)}(g) := g + \varepsilon\{g, f\} + O(\varepsilon^2)$$

- $\Phi_f^{(e)}$ is **generator** of map from \mathcal{P} to \mathcal{P} :

$$\Phi_f^{(e)}(q_i) = q_i + \varepsilon \frac{\partial f}{\partial p_i} + O(\varepsilon^2)$$

$$\Phi_f^{(e)}(p_i) = p_i - \varepsilon \frac{\partial f}{\partial q_i} + O(\varepsilon^2)$$

- Noether charge $Q = \left(\sum_{i=1}^n a_i p_i\right) - F$ generates symmetry transformation via Hamiltonian flow:

$$\Phi_Q^{(e)}(q_i) = q_i + \varepsilon\{q_i, Q\} + O(\varepsilon^2) = q_i + \varepsilon a_i + O(\varepsilon^2)$$

- **Hamiltonian** gives **energy**:

$$H = \left(\sum_{i=1}^n p_i \dot{q}_i\right) - L$$

- **Hamilton's equations of motion:**

$$\dot{q}_i = \{q_i, H\} = \frac{\partial H}{\partial p_i}, \quad \dot{p}_i = \{p_i, H\} = -\frac{\partial H}{\partial q_i}$$

- Time evolution of $f(\underline{q}, \underline{p})$ generated by H :

$$\frac{df}{dt} = \{f, H\}$$

In f depends explicitly on time,

$$\frac{df}{dt} = \frac{\partial f}{\partial t} + \{f, H\}$$

- **Relation between Hamiltonian and Lagrangian:**

$$\left. \frac{\partial H(\underline{q}, \underline{p}, t)}{\partial t} \right|_{\underline{q}, \underline{p}} = - \left. \frac{\partial L(\underline{q}, \underline{\dot{q}}, t)}{\partial t} \right|_{\underline{q}, \underline{\dot{q}}}$$

- If function Q doesn't depend explicitly on time, $\{H, Q\} = 0$ so Hamiltonian left invariant by transformation generated by Q :

$$\Phi_Q(H) = H + \varepsilon\{Q, H\} + O(\varepsilon^2) = H + O(\varepsilon^2)$$

6. Wave function and probabilities

- **Wave function:** continuous, complex function of position x and time t : $\psi(x, t)$.
- **Probability density to find particle at time t and position x :** $P(x, t) = |\psi(x, t)|^2$, with

$$\int_{-\infty}^{\infty} P(x, t) dx = 1$$

If this integral exists, ψ is **square-normalisable**. If integral equal to 1, ψ is **normalised**. Probability of finding particle in interval (a, b) is

$$\int_a^b P(x, t) dx$$

- **Expectation value of $f(x)$:**

$$\langle f(x) \rangle = \int_{-\infty}^{\infty} f(x) P(x, t) dx$$

- **Uncertainty in position:** $\Delta x = \sqrt{\langle x^2 \rangle - \langle x \rangle^2}$
- **Infinite potential well in $0 < x < L$:**

$$V(x) = \begin{cases} 0 & \text{if } 0 < x < L \\ \infty & \text{otherwise} \end{cases}$$

Wave function vanishes in regions $x \leq 0$ and $x \geq L$. Eigenfunctions for this potential are

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

- **Wave function collapse:** if position is measured to be x_0 , wave function becomes very localised around at x_0 , and measurement immediately afterwards will also yield x_0 .
- $\langle x \rangle$ is not average of repeated measurements of same particle, but average of measurements of many particles with same wave function.

7. Momentum and Planck's constant

- **Position operator:**

$$\hat{x} = x$$

- **Momentum operator:**

$$\hat{p} = -i\hbar \frac{\partial}{\partial x}$$

where \hbar is **reduced Planck constant**.

- **Commutator:**

$$[\hat{x}, \hat{p}] := \hat{x}\hat{p} - \hat{p}\hat{x} = i\hbar$$

- **Expectation value of momentum for wave function ψ :**

$$\langle \psi \rangle = \int_{-\infty}^{\infty} \overline{\psi(x, t)} \hat{p} \psi(x, t) dx = -i\hbar \int_{-\infty}^{\infty} \overline{\psi(x, t)} \frac{\partial}{\partial x} \psi(x, t) dx$$

- **Expection value of function of momentum:**

$$\langle f(p) \rangle = \int_{-\infty}^{\infty} \overline{\psi(x, t)} f(\hat{p}) \psi(x, t) dx$$

- **Momentum uncertainty:**

$$\Delta p = \sqrt{\langle p^2 \rangle - \langle p \rangle^2}$$

- **Heisenberg's uncertainty principle:** for any normalised wave function,

$$\Delta x \Delta p \geq \frac{\hbar}{2}$$

8. Schrodinger's equation

- **Hamiltonian operator:**

$$\hat{H} = \frac{\hat{p}^2}{2m} + V(x) = -\frac{\hbar^2}{2m} \frac{\partial^2}{\partial x^2} + V(x)$$

Corresponds to measurements of energy.

- **Schrodinger's equation:**

$$i\hbar \frac{\partial \psi(x, t)}{\partial t} = \hat{H} \psi(x, t)$$

9. The Hilbert space

- **Hermitian inner product on vector space V :** map $\langle \cdot, \cdot \rangle : V \times V \rightarrow \mathbb{C}$ satisfying:

- $\langle v, w \rangle = \overline{\langle w, v \rangle}$.
- $\langle v, a_1 w_1 + a_2 w_2 \rangle = a_1 \langle v, w_1 \rangle + a_2 \langle v, w_2 \rangle$.
- $\langle a_1 v_1 + a_2 v_2, w \rangle = \overline{a_1} \langle v_1, w \rangle + \overline{a_2} \langle v_2, w \rangle$
- $\langle v, v \rangle \geq 0$ for all v and $\langle v, v \rangle = 0 \iff v = 0$.

- Set of continuous square-integrable wave functions forms complex vector space. So $a_1 \psi_1 + a_2 \psi_2$ is also square-integrable.

- **Hermitian inner product of two wave functions:**

$$\langle \psi_1, \psi_2 \rangle = \int_{-\infty}^{\infty} \overline{\psi_1(x)} \psi_2(x) dx$$

- If $\{\varphi_n(x)\}$ is orthonormal basis so $\langle \varphi_m, \varphi_n \rangle = \delta_{mn}$, then any vector can be expressed

$$\psi(x) = \sum_n c_n \varphi_n(x)$$

where $c_m = \langle \varphi_m, \psi \rangle$. Hermitian product is then

$$\langle \psi_1, \psi_2 \rangle = \sum_i \overline{\psi_1(x)} \psi_2(x) = \sum_n \bar{c}_{1,n} c_{2,n}$$

So squared norm of ψ is $|\psi|^2 = \langle \psi, \psi \rangle = \sum_n |c_n|^2$.

10. Hermitian operators

- For vector space V , **linear operator** is map $A : V \rightarrow V$ with

$$A(a_1 v_1 + a_2 v_2) = a_1 (A v_1) + a_2 (A v_2)$$

- Any linear combination or composition of linear operators is linear operator.
- Matrix elements of linear operator** for orthonormal basis $\{e_j\}$: $A_{ij} = \langle e_i, A e_j \rangle$.
- Adjoint** A^\dagger : $\langle v_1, A v_2 \rangle = \langle A^\dagger v_1, v_2 \rangle$. Adjoint has matrix elements which are conjugate of transpose of original matrix.
- Properties of adjoint:
 - $(a_1 A_1 + a_2 A_2)^\dagger = \overline{a_1} A_1^\dagger + \overline{a_2} A_2^\dagger$.
 - $(A_1 A_2)^\dagger = A_2^\dagger A_1^\dagger$.
- Hermitian operator**: linear operator that is equal to adjoint. Matrix is Hermitian: $A_{ij} = \overline{A_{ji}}$.
- Position and momentum operators Hermitian, w.r.t. orthonormal basis of wave functions $\{\varphi_n(x)\}$.

11. The spectrum of a Hermitian operator

- Wave function ψ_a is **eigenfunction** of Hermitian differential operator A with **eigenvalue** a if $A \psi_a(x) = a \psi_a(x)$.
- Expectation value of Hermitian operator:

$$\langle A \rangle = \langle \psi, A \psi \rangle = \int_{-\infty}^{\infty} \overline{\psi(x)} A \psi(x) dx$$

- .
- If ψ_a is eigenfunction, $\langle A \rangle = a$ and $\langle A^n \rangle = a^n$. So uncertainty $\Delta A = 0$.
- Let A Hermitian operator.
 - Eigenvalues are real and
 - ψ_1, ψ_2 eigenfunctions of A with distinct eigenvalues are orthogonal.
- If A has discrete spectrum, can choose orthonormal basis of eigenfunctions $\{\varphi_n(x)\}$ with eigenvalues a_n . Then any wave function can be written as $\psi(x) = \sum_n c_n \varphi_n(x)$ where $c_n = \langle \varphi_n, \psi \rangle$. Can interpret $|c_n|^2$ as probability of measurement of A yielding a_n .
- Dirac delta function**:

$$\delta(a) = \begin{cases} 0 & \text{if } a \neq 0 \\ \infty & \text{if } a = 0 \end{cases}$$

with $\int_{-\infty}^{\infty} \delta(a) da = 1$ and

$$\int_{-\infty}^{\infty} \delta(a - a') f(a') da' = f(a)$$

- Limit definition of Dirac delta function**: limit as $\varepsilon \rightarrow 0^+$ of

$$\delta_\varepsilon(a) = \frac{1}{\varepsilon\sqrt{\pi}} e^{-a^2/\varepsilon^2}$$

- **Delta function is Fourier transform of 1:**

$$\delta(a) = \frac{1}{2\pi} \int_{-\infty}^{\infty} e^{iaa'} da'$$

- If A has continuous spectrum (eigenvalues $a \in \mathbb{R}$) then can choose basis of eigenfunctions $\varphi_a(x)$ with $\langle \varphi_a, \varphi_{a'} \rangle = \delta(a - a')$. Can uniquely expand wave function

$$\psi(x) = \int_{-\infty}^{\infty} c(a) \varphi_a(x) da$$

where $c(a) = \langle \varphi_a, \psi \rangle$. Norm of wave function is

$$\langle \psi, \psi \rangle = \int_{-\infty}^{\infty} |c(a)|^2 da$$

For normalised wave function,

$$\int_{-\infty}^{\infty} |c(a)|^2 da = 1$$

so treat $|c(a)|^2$ as probability distribution for measurements of A .

12. Postulates of quantum mechanics

- **Postulates of quantum mechanics:**

- Particle described by normalised wave function $\psi(x)$.
- Measurable quantities represented by Hermitian operators $A(x, p)$, constructed from polynomial/real analytic functions of position and momentum operators:

$$\hat{x} = x,$$

$$\hat{p} = -i\hbar \frac{\partial}{\partial x}$$

- Possible outcomes of measurement of A are given by its eigenvalues a . If spectrum discrete, $\{a_j\}$, then choose eigenfunction basis $\varphi_j(x)$ with $\langle \varphi_i, \varphi_j \rangle = \delta_{ij}$. Then probability of finding measurement as eigenvalue a_j is $|\langle \varphi_j, \psi \rangle|^2$. If spectrum continuous, $a \in \mathbb{R}$, choose eigenfunctions $\varphi_a(x)$ with $\langle \varphi_a, \varphi_{a'} \rangle = \delta(a - a')$, then probability of finding measurement as eigenvalue a is $|\langle \varphi_a, \psi \rangle|^2$.
- If measurement of A yields eigenvalue a_j (or a), wave function immediately afterwards is $\varphi_j(x)$ (or $\varphi_a(x)$). **Note:** in continuous case, wave function immediately afterwards not square-normalisable.
- If no measurements made, ψ evolves in time according to Schrodinger equation:

$$i\hbar \frac{\partial \psi(x, t)}{\partial t} = \hat{H} \psi(x, t)$$

- For discrete spectrum, expectation value of A is

$$\langle A \rangle = \sum_j a_j P_j$$

for eigenvalues a_j , $P_j = |c_j|^2$ is probability of measurement being a_j .

- For continuous spectrum, expectation value of A is

$$\langle A \rangle = \int_{-\infty}^{\infty} a P(a) da$$

where $P(a) = |c(a)|^2$ is probability distribution.

13. Commutators and uncertainty principle

- **Commutator** of operators A, B :

$$[A, B] = AB - BA$$

- Properties of commutator:
 - **Anti-symmetry**: $[A, B] = -[B, A]$.
 - **Linearity**: $[a_1 A_1 + a_2 A_2, B] = a_1 [A_1, B] + a_2 [A_2, B]$.
 - $[A, BC] = B[A, C] + [A, B]C$.
 - **Jacobi identity**: $[A, [B, C]] + [B, [C, A]] + [C, [A, B]] = 0$.
- If $[A, B] = 0$, possible to find orthonormal basis of wave functions which are eigenfunctions of A and B .
- A, B **compatible** if $[A, B] = 0$.
- **Generalised uncertainty principle**: for any square-normalisable wave function,

$$\Delta A \Delta B \geq \frac{1}{2} |\langle [A, B] \rangle|$$

- **Anti-commutator**: $\{A, B\} = AB + BA$.

14. Energy revisited

- Eigenfunctions of Hamiltonian are **bound states** if classical solution is bounded in space.
- Let $V(x) \geq V_0$ for all $x \in \mathbb{R}$. Then if wave function normalised, $\langle H \rangle > V_0$.
- If $\psi(x)$ is normalised eigenfunction of H with eigenvalue E , then $E > V_0$.
- **Zero-point energy**: smallest eigenvalue $E > V_0$.
- Spectrum of Hamiltonian is non-degenerate.

15. Stationary states

- Solution to Schrodinger's equation is $\psi(x, t) = \varphi(x)e^{-iEt/\hbar}$ where $\varphi(x)$ is eigenfunction of Hamiltonian with eigenvalue E . This solution is **stationary wave function**.
- **Full solution to Schrodinger's equation**:

$$\psi(x, t) = \sum_j c_j \varphi_j(x) e^{-iE_j t/\hbar}$$

where $\{\varphi_j(x)\}$ is orthonormal basis of Hamiltonian eigenfunctions with eigenvalues E_j , c_j are coefficients of initial wave function expansion:

$$\psi(x, 0) = \sum_j c_j \varphi_j(x)$$

Probability of energy measurement being E_j is $P_j = |\langle \varphi_j, \psi \rangle|^2 = |c_j|^2$.

16. Case study: the free particle

- If $V(x) = 0$, eigenfunction of \hat{p} is eigenfunction of \hat{H} .

17. Two particle systems

- For two particles in one dimension, wave function is $\psi(x_1, x_2)$, probability density is $P(x_1, x_2) = |\psi(x_1, x_2)|^2$: probability of finding particle one in (a, b) and particle two in (c, d) is

$$\int_a^b \int_c^d P(x_1, x_2) dx_1 dx_2$$

- Probability of finding particle one in (a, b) is

$$P(x_1) = \int_a^b P(x_1, x_2) dx_2$$

(similarly for particle two).

- If both positions measured as \tilde{x}_1, \tilde{x}_2 , wave function collapses to product of position eigenfunctions:

$$\psi_{\text{before}}(x_1, x_2) \rightarrow \psi_{\text{after}}(x_1, x_2) \propto \delta(x_1 - \tilde{x}_1) \delta(x_2 - \tilde{x}_2)$$

- If only particle one measured,

$$\psi_{\text{before}}(x_1, x_2) \rightarrow \delta(x_1 - \tilde{x}_1) \psi_{\text{before}}(\tilde{x}_1, x_2)$$

- Hamiltonian for two particles with zero potential:

$$\hat{H} = -\frac{\hbar^2}{2m} \frac{\partial}{\partial x_1^2} - \frac{\hbar^2}{2m} \frac{\partial}{\partial x_2^2}$$

Eigenfunctions are product are single-particle eigenfunctions:

$$\varphi(x_1, x_2) = \frac{2}{L} \sin\left(\frac{n\pi x_1}{L}\right) \sin\left(\frac{m\pi x_2}{L}\right)$$

Eigenvalues are sum of eigenvalues of single-particle Hamiltonians.

- Wave function separable if can be written as product of function of x_1 and function of x_2 .
- **Entangled states:** when measurement of one particle affects subsequent measurement of other particle. Occurs for non-separable wave functions.

18. Simple harmonic oscillator

- Simple harmonic oscillator potential:

$$V(x) = \frac{1}{2}m\omega^2 x^2$$

where ω is angular frequency.

- If V has minimum at $x = x_0$ and $|x - x_0|$ small, $m\omega^2 \approx \frac{1}{2}V''(x_0)$ by Taylor expansion of $V(x)$ around x_0 .
- Energy spectrum of Hamiltonian for simple harmonic oscillator is $E_n = \hbar\omega(n + \frac{1}{2})$