

Notes of Quantum Mechanics

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Preface

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Chapter 1

One-dimensional harmonic oscillator

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1.1 Introduction

1.1.1 Importance of the harmonic oscillator in physics

The simplest example is a particle of mass m moving in a potential which depends only on x and has the form

$$V(x) = \frac{1}{2}kx^2, \quad k > 0.$$

The particle is attracted towards the $x = 0$ by a restoring force:

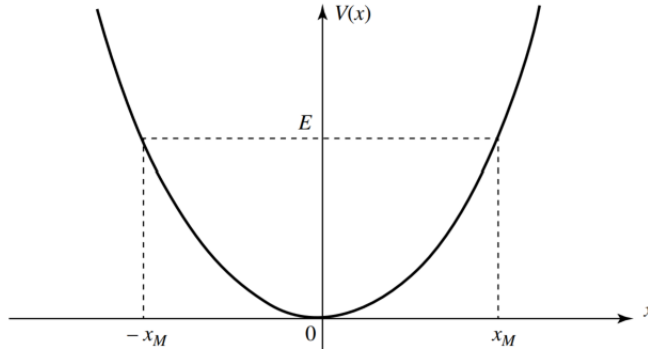


Figure 1.1 Potential energy $V(x)$ of a 1D harmonic oscillator.

$$F_x = \frac{dV}{dx} = -kx.$$

In classical mechanics, the motion of the particle is a sinusoidal oscillation about $x = 0$ with angular frequency $\omega = \sqrt{k/m}$.

Various systems are governed by the harmonic oscillator equations

Whenever one studies the behavior of a system in the neighborhood of a stable equilibrium position, one arrives at equations which, in the limit of small oscillations, are those of a harmonic oscillator.

1.1.2 The harmonic oscillator in classical mechanics

The motion of the particle is governed by the dynamics equation

$$m \frac{d^2x}{dt^2} = -\frac{dV}{dx} = -kx \longrightarrow x = x_M \cos(\omega t - \varphi). \quad (1.1)$$

The kinetic energy of the particle is

$$T = \frac{1}{2}m \left(\frac{dx}{dt} \right)^2 = \frac{p^2}{2m}, \quad (1.2)$$

where $p = mv$ is the momentum of the particle. The total energy is

$$E = T + V = \frac{p^2}{2m} + \frac{1}{2}m\omega^2 x^2 = \frac{1}{2}m\omega^2 x_M^2.$$

- The potential can be expanded in Taylor's series around x_0 :

$$V(x) = \underbrace{V(x_0)}_a + V'(x_0)(x - x_0) + \underbrace{\frac{1}{2!}V^{(2)}(x_0)(x - x_0)^2}_b + \underbrace{\frac{1}{3!}V^{(3)}(x_0)(x - x_0)^3}_c + \dots$$

The force derived from the potential in the neighborhood of x_0 is

$$F_x = -\frac{dV}{dx} = -2b(x - x_0) - 3c(x - x_0)^2 + \dots \quad (1.3)$$

The point $x = x_0$ is a stable equilibrium for the particle: $F_x(x_0) = 0$. In addition, if the amplitude of the motion of the particle about x_0 is sufficiently small, we can keep with the linear term only and we have a harmonic oscillator since the dynamics equation can be approximated by

$$m \frac{d^2x}{dt^2} \approx -2b(x - x_0).$$

For higher energies E , the particle will be in period but not sinusoidal motion (as signal in Fourier series) between the limits x_1 and x_2 . We then say that we are dealing with an **anharmonic oscillator**.

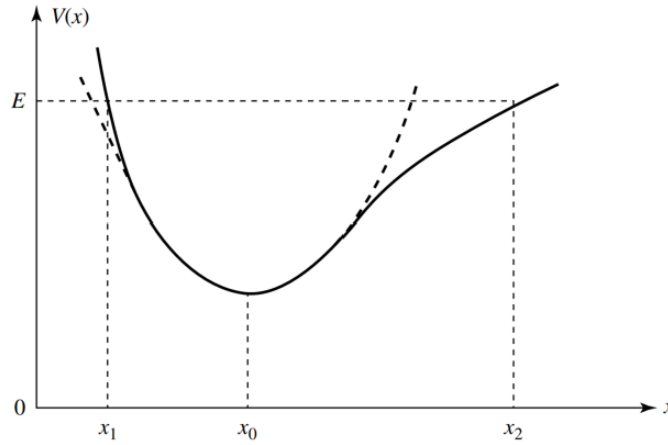


Figure 1.2 Any potential can be approximated by a parabolic potential. In $V(x)$, a classical particle of energy E oscillates between x_1 and x_2 .

1.1.3 General properties of the quantum mechanical Hamiltonian

In QM, the classical quantities x and p are replaced respectively by the observables X and P , which satisfy

$$[X, P] = i\hbar.$$

It is then easy to obtain the Hamiltonian operator of the system from the total energy

$$H = \frac{P^2}{2m} + \frac{1}{2}m\omega^2 X^2.$$

Since H is time-independent (conservative system), the quantum mechanical study of the harmonic oscillator reduces to the solution of the eigenequation:

$$H|\varphi\rangle = E|\varphi\rangle$$

which is written, in the $\{|x\rangle\}$ representation

$$\left[-\frac{\hbar^2}{2m} \frac{d^2}{dx^2} + \frac{1}{2} m \omega^2 x^2 \right] \varphi(x) = E \varphi(x).$$

Let us indicate some properties of the potential function:

- **The eigenvalues of the Hamiltonian are positive.** If $V(x)$ has a lower bound, the eigenvalues E of H are greater than the minimum of $V(x)$:

$$V(x) \leq V_m \quad \text{requires} \quad E > V_m.$$

We have chosen for the harmonic oscillator that $V_m = 0$.

- **The eigenfunctions of H have a definite parity** due to that $V(-x) = V(x)$ is an even function. We shall see that the eigenvalues of H are not degenerate; the wave functions associated with the stationary states are necessarily either even or odd.
- **The energy spectrum is discrete.**

1.2 Eigenvalues of the Hamiltonian

1.2.1 Notation

It is easy to see that the observables \hat{X} and \hat{P}

$$\text{dimensionless observables} \quad \hat{X} = \sqrt{\frac{m\omega}{\hbar}} X, \quad \hat{P} = \frac{1}{\sqrt{m\hbar\omega}} P$$

are dimensionless. With these new operators, the canonical commutation is

$$\text{Canonical commutation} \quad [\hat{X}, \hat{P}] = i \quad (1.4)$$

and the Hamiltonian can be put in the form

$$H = \hbar\omega \hat{H}, \quad \text{with} \quad \hat{H} = \frac{1}{2}(\hat{X}^2 + \hat{P}^2). \quad (1.5)$$

In consequence, we seek the solutions of the following eigenequation

$$\hat{H}|\varphi_\nu^i\rangle = \epsilon_\nu |\varphi_\nu^i\rangle,$$

where the operator \hat{H} and the eigenvalues ϵ_ν are **dimensionless**.

If \hat{X} and \hat{P} were numbers and not operators, we could write the sum $\hat{X}^2 + \hat{P}^2$ appearing in the definition of \hat{H} in the form of a product $(\hat{X} - i\hat{P})(\hat{X} + i\hat{P})$. However, the introduction of operators proportional to $\hat{H} \pm i\hat{P}$ enables us to simplify considerably our search for eigenvalues and eigenvectors of \hat{H} . We therefore set

$$\begin{aligned} a &= \frac{1}{\sqrt{2}}(\hat{X} + i\hat{P}) & \hat{X} &= \frac{1}{\sqrt{2}}(a^\dagger + a) \\ a^\dagger &= \frac{1}{\sqrt{2}}(\hat{X} - i\hat{P}) & \hat{P} &= \frac{i}{\sqrt{2}}(a^\dagger - a) \end{aligned} \quad \Longleftrightarrow \quad (1.6)$$

The commutator of a and a^\dagger is

$$[a, a^\dagger] = \frac{1}{2}[\hat{X} + i\hat{P}, \hat{X} - i\hat{P}] = \frac{i}{2}[\hat{P}, \hat{X}] - \frac{i}{2}[\hat{X}, \hat{P}] = 1 \longrightarrow [a, a^\dagger] = 1. \quad (1.7)$$

If we do aa^\dagger we obtain

$$a^\dagger a = \frac{1}{2}(\hat{X} - i\hat{P})(\hat{X} + i\hat{P}) = \frac{1}{2}(\hat{X}^2 + \hat{P}^2 + i\hat{X}\hat{P} - i\hat{P}\hat{X}) = \frac{1}{2}(\hat{X}^2 + \hat{P}^2 - 1).$$

Comparing with \hat{H} we see that

$$\hat{H} = a^\dagger a + \frac{1}{2} = aa^\dagger - \frac{1}{2}.$$

We see that we cannot put \hat{H} in a product of linear terms, due to the non-commutativity of \hat{X} and \hat{P} (1/2 term).

We introduce another operator:

$$\text{Operator } N \quad N = a^\dagger a. \quad (1.8)$$

This operator is Hermitian

$$N^\dagger = a^\dagger (a^\dagger)^\dagger = a^\dagger a = N. \quad (1.9)$$

And its relation with \hat{H} is

$$\hat{H} = N + \frac{1}{2} \quad (1.10)$$

so that the eigenvectors of \hat{H} are eigenvectors of N , and viceversa. The commutators with a and a^\dagger are:

$$[N, a] = [a^\dagger a, a] = a^\dagger [a, a] + [a^\dagger, a]a = -a \longrightarrow [N, a] = -a \quad (1.11)$$

$$[N, a^\dagger] = [a^\dagger a, a^\dagger] = a^\dagger [a, a^\dagger] + [a^\dagger, a^\dagger]a = a^\dagger \longrightarrow [N, a^\dagger] = a^\dagger. \quad (1.12)$$

The study of the harmonic oscillator is based on these operators a , a^\dagger , and N . The eigenequation for N is

$$\text{Eigenequation of } N \quad N|\varphi_\nu^i\rangle = \nu|\varphi_\nu^i\rangle. \quad (1.13)$$

When this is solved, we know that the eigenvector $|\varphi_\nu^i\rangle$ of N is also an eigenvector of H with the eigenvalue $E_\nu = (\nu + 1/2)\hbar\omega$:

$$H|\varphi_\nu^i\rangle = (\nu + 1/2)\hbar\omega|\varphi_\nu^i\rangle. \quad (1.14)$$

The solution of the eigenequation of N will be based on the commutation relation $[a, a^\dagger] = 1$.

1.2.2 Determination of the spectrum

Lemmas

- **Properties of the eigenvalues of N** The eigenvalues ν of the operator N are positive or zero. We can see this by looking the square of the norm of the vector $a|\varphi_n^i\rangle$

$$\|a|\varphi_\nu^i\rangle\|^2 = \langle\varphi_\nu^i|a^\dagger a|\varphi_\nu^i\rangle = \langle\varphi_\nu^i|N|\varphi_\nu^i\rangle = \nu\langle\varphi_\nu^i|\varphi_\nu^i\rangle \geq 0 \implies \nu \geq 0.$$

- **Properties of the vector $a|\varphi_\nu^i\rangle$**

- $\nu = 0 \implies a|\varphi_{\nu=0}^i\rangle = 0$. If $\nu = 0$ is an eigenvalue of N , all eigenvectors $|\varphi_0^i\rangle$ associated with this eigenvalue satisfy the relation

$$a|\varphi_0^i\rangle = 0. \quad (1.15)$$

Any vector which satisfy this relation is therefore an eigenvector of N with the eigenvalue $\nu = 0$.

- $\nu > 0 \implies a|\varphi_\nu^i\rangle$ is a non-zero eigenvector of N with eigenvalue $\nu - 1$.

$$\begin{aligned} [N, a]|\varphi_\nu^i\rangle &= -a|\varphi_\nu^i\rangle \\ Na|\varphi_\nu^i\rangle &= aN|\varphi_\nu^i\rangle - a|\varphi_\nu^i\rangle \implies N[a|\varphi_\nu^i\rangle] = (\nu - 1)[a|\varphi_\nu^i\rangle] \\ N[a|\varphi_\nu^i\rangle] &= a\nu|\varphi_\nu^i\rangle - a|\varphi_\nu^i\rangle \end{aligned}$$

- **Properties of the vector $a^\dagger|\varphi_\nu^i\rangle$**

- $a^\dagger|\varphi_\nu^i\rangle$ is always non-zero. We study it with the square of the norm:

$$\|a^\dagger|\varphi_\nu^i\rangle\|^2 = \langle\varphi_\nu^i|aa^\dagger|\varphi_\nu^i\rangle = \langle\varphi_\nu^i|(N + 1)|\varphi_\nu^i\rangle = (\nu + 1)\langle\varphi_\nu^i|\varphi_\nu^i\rangle.$$

As $\nu \geq 0$ by lemma 1, the ket $a^\dagger|\varphi_\nu^i\rangle$ always has non-zero norm and, consequently, is never zero.

- $a^\dagger|\varphi_\nu^i\rangle$ is an eigenvector of N with eigenvalue $N + 1$. We do it analogously to lemma IIb):

$$\begin{aligned} [N, a^\dagger]|\varphi_\nu^i\rangle &= a^\dagger|\varphi_\nu^i\rangle \\ Na^\dagger|\varphi_\nu^i\rangle &= a^\dagger N|\varphi_\nu^i\rangle + a^\dagger|\varphi_\nu^i\rangle \implies N[a^\dagger|\varphi_\nu^i\rangle] = (\nu + 1)[a^\dagger|\varphi_\nu^i\rangle] \\ N[a^\dagger|\varphi_\nu^i\rangle] &= \nu a^\dagger|\varphi_\nu^i\rangle + a^\dagger|\varphi_\nu^i\rangle \end{aligned}$$

The spectrum of N is composed of non-negative integers

If ν is non-integral, we can therefore construct a non-zero eigenvector of N with a strictly negative eigenvalue. Since this is impossible by lemma 1, the hypothesis of non-integral ν must be rejected.

ν can only be a non-negative integer.

We conclude that the eigenvalues of H are of the form

$$\text{Eigenvalue of } H \quad E_n = \left(n + \frac{1}{2}\right) \hbar\omega, \quad n \in \mathbb{N}_0^+. \quad (1.16)$$

In QM, the energy of the harmonic oscillator is **quantized**. The smallest value (ground state) is $\hbar\omega/2$.

Interpretation of the a and a^\dagger operators

We have seen that, given $|\varphi_n^i\rangle$ with eigenvalue E_n , application of a gives an eigenvector associated with E_{n-1} while application of a^\dagger yields the energy E_{n+1} .

That's why a^\dagger is said to be a **creation operator** and a an **annihilation operator**; their action on an eigenvector of N makes an energy quantum $\hbar\omega$ appear or disappear.

1.2.3 Degeneracy of the eigenvalues

The ground state is non-degenerate

The eigenstates of H associated with $E_0 = \hbar\omega/2$ (or eigenvector of N associated with $n = 0$), according to lemma II, must all satisfy the equation

$$a|\varphi_0^i\rangle = 0.$$

To find the degeneracy of the E_0 level, all we must do is see how many kets satisfy the above. We can write the above equation using the definition of \hat{X} , \hat{P} and a in terms of them, in the form

$$\frac{1}{\sqrt{2}} \left[\sqrt{\frac{m\omega}{\hbar}} X + \frac{i}{\sqrt{m\hbar\omega}} P \right] |\varphi_0^i\rangle = 0.$$

In the $\{|x\rangle\}$ representation, this relation becomes

$$\left(\frac{m\omega}{\hbar} x + \frac{d}{dx} \right) \varphi_0^i(x) = 0, \quad \text{where} \quad \varphi_0^i(x) = \langle x | \varphi_0^i \rangle.$$

Therefore we must solve a first-order differential equation, whose solution is

$$\varphi_0^i(x) = c e^{-\frac{1}{2} \frac{m\omega}{\hbar} x^2} \quad (1.17)$$

The various solutions of the ODE are all proportional to each other. Consequently, there exists only one ket $|\varphi_0\rangle$ that satisfies the initial equation: the ground state $E_0 = \hbar\omega/2$ is not degenerate.

All the states are non-degenerate

We use recurrence to show that all other states are also non-degenerate. We need to prove that if E_n is non-degenerate, the level E_{n+1} is not either.

Let's assume there exists only one vector $|\varphi_n\rangle$ such that

$$N|\varphi_n\rangle = n|\varphi_n\rangle.$$

Then consider an eigenvector $|\varphi_{n+1}^i\rangle$ corresponding to the eigenvalue $n+1$

$$N|\varphi_{n+1}^i\rangle = (n+1)|\varphi_{n+1}^i\rangle.$$

We know that the ket $a|\varphi_{n+1}^i\rangle$ is not zero and that it is an eigenvector of N with eigenvalue n . Since this ket is not degenerate by hypothesis, there exists a number c^i such that

$$a|\varphi_{n+1}^i\rangle = c^i|\varphi_n\rangle / a^\dagger \longrightarrow a^\dagger a|\varphi_{n+1}^i\rangle = N|\varphi_{n+1}^i\rangle = (n+1)|\varphi_{n+1}^i\rangle = c^i a^\dagger |\varphi_n\rangle.$$

We have,

$$|\varphi_{n+1}^i\rangle = \frac{c^i}{n+1} a^\dagger |\varphi_n\rangle.$$

We see that all kets $|\varphi_{n+1}^i\rangle$ associated with the eigenvalue $n + 1$ are proportional to $a^\dagger|\varphi_n\rangle$. They are proportional to each other: the eigenvalue $n + 1$ is not degenerate.

Since the eigenvalue $n = 0$ is not degenerate, the eigenvalue $n = 1$ is not either, nor is $n = 2$, etc.: all the eigenvalues of N and, consequently, all those of H , are non-degenerate. Now, we can just write $|\varphi_n\rangle$ for the eigenvector of H associated with E_n .

1.3 Eigenstates of the Hamiltonian

1.3.1 The $\{\varphi_n\}$ representation

Since none of the eigenvalues of N (H) is degenerate, N (H) alone constitutes a CSCO in \mathcal{E}_c .

The basis vectors in terms of $|\psi_0\rangle$

We assume that the vector $|\varphi_0\rangle$ which satisfies $a|\varphi_0\rangle = 0$, is normalized. According to lemma III, the vector $|\varphi_1\rangle$ is proportional to $a^\dagger|\varphi_0\rangle$ in the form

$$|\varphi_1\rangle = c_1 a^\dagger |\varphi_0\rangle.$$

We shall determine c_1 by requiring $|\varphi_1\rangle$ to be normalized and choosing the phase of $|\varphi_1\rangle$ such that c_1 is real and positive. The square of the norm of $|\varphi_1\rangle$ is

$$\langle\varphi_1|\varphi_1\rangle = |c_1|^2 \langle\varphi_0|aa^\dagger|\varphi_0\rangle = |c_1|^2 \langle\varphi_0|(a^\dagger a + 1)|\varphi_0\rangle = |c_1|^2 [\underbrace{\langle\varphi_0|N|\varphi_0\rangle}_{0\langle\varphi_0|\varphi_0\rangle} + \langle\varphi_0|\varphi_0\rangle] = |c_1|^2.$$

We find that $c_1 = 1$:

$$\langle\varphi_1|\varphi_1\rangle = |c_1|^2 = 1 \implies |\varphi_1\rangle = a^\dagger|\varphi_0\rangle. \quad (1.18)$$

We can do the same to construct $|\varphi_2\rangle$ from $|\varphi_1\rangle$ and get c_2 and so on. In general, if we know $|\varphi_{n-1}\rangle$ (normalized), then the normalized vector $|\varphi_n\rangle$ is written

$$|\varphi_n\rangle = c_n a^\dagger |\varphi_{n-1}\rangle, \quad \text{so that} \quad c_n = \frac{1}{\sqrt{n}}.$$

In fact, we can express all $|\varphi_n\rangle$ in terms of $|\varphi_0\rangle$ by recursion:

$$|\varphi_n\rangle = \frac{1}{\sqrt{n!}} (a^\dagger)^n |\varphi_0\rangle. \quad (1.19)$$

Orthonormalization and closure relations

Action of the various operators

1.3.2 Wave functions associated with the stationary states

1.4 Discussion

1.4.1 Mean values and rms deviations of X and P in a state $|\varphi_n\rangle$

1.4.2 Properties of the ground state

1.4.3 Time evolution of the mean values

Formula sheet

1.4.4 Useful formulas

Closure relation (discrete)	$\sum_k \sum_{i=1}^{g_k} v_k^i\rangle \langle v_k^i = \mathbb{1}$	Closure relation (continuous)	$\int_{\beta} d\beta \omega_{\beta}\rangle \langle \omega_{\beta} = \mathbb{1}$
Glauber Formula	$e^{A+B} = e^A e^B e^{-\frac{1}{2}[A,B]}$	Generalized uncertainty relation	$\Delta A \Delta B \geq \frac{1}{2} \langle [A, B] \rangle $
Function of an operator	$F(A) = \sum_{n=0}^{\infty} f_n (A - a)^n$		$\Delta Q = \sqrt{\langle Q^2 \rangle - \langle Q \rangle^2}$
Eigenequation of $F(A)$	$F(A) \psi\rangle = F(\lambda) \psi\rangle$		
Transformation $\{u\} \rightarrow \{v\}$	$\mathbb{M}_{jk} = \langle u_j v_k \rangle$	$ \psi\rangle_{\{u\}} = \mathbb{M} \psi\rangle_{\{v\}}$ $A_{\{u\}} = \mathbb{M} A_{\{v\}} \mathbb{M}^{\dagger}$	$ \psi\rangle_{\{v\}} = \mathbb{M}^{\dagger} \psi\rangle_{\{u\}}$ $A_{\{v\}} = \mathbb{M}^{\dagger} A_{\{u\}} \mathbb{M}$

1.4.5 Basis

Quantity	Discrete basis (sum over j, k)	Continuous basis (integrate over β, β')
$\mathbb{1}$	$= \sum v_k\rangle \langle v_k $	$= \int d\beta \omega_{\beta}\rangle \langle \omega_{\beta} $
$ \psi\rangle = \mathbb{1} \psi\rangle$	$= \sum v_k\rangle \langle v_k \psi\rangle$	$= \int d\beta \omega_{\beta}\rangle \langle \omega_{\beta} \psi\rangle$
$\langle \varphi = \langle \varphi \mathbb{1}$	$= \sum \langle \varphi v_k\rangle \langle v_k $	$= \int d\beta \langle \varphi \omega_{\beta}\rangle \langle \omega_{\beta} $
$A = \mathbb{1}A\mathbb{1}$	$= \sum \sum v_j\rangle \langle v_j A v_k\rangle \langle v_k $	$= \iint d\beta d\beta' \omega_{\beta}\rangle \langle \omega_{\beta} A \omega_{\beta'}\rangle \langle \omega_{\beta'} $

Quantity	X representation	P_x representation
X	x	$i\hbar \partial/\partial p$
P_x	$-i\hbar \partial/\partial x$	p
$ x'\rangle$	$\langle x x'\rangle = \delta(x - x')$	$\langle p x'\rangle = \frac{1}{\sqrt{2\pi\hbar}} \exp(-ix'p/\hbar)$
$ p'\rangle$	$\langle x p'\rangle = \frac{1}{\sqrt{2\pi\hbar}} \exp(ixp'/\hbar)$	$\langle p p'\rangle = \delta(p - p')$
$ \psi\rangle$	$\langle x \psi\rangle = \psi(x)$	$\langle p \psi\rangle = \tilde{\psi}(p)$

Fourier transforms for 3D wavefunctions

$\tilde{\psi}(\mathbf{p}) = \mathcal{F}[\psi(\mathbf{r})] = \left(\frac{1}{2\pi\hbar}\right)^{3/2} \int_{-\infty}^{\infty} d^3\mathbf{r} e^{-i\mathbf{r}\cdot\mathbf{p}/\hbar} \psi(\mathbf{r})$	$\psi(\mathbf{r}) = \mathcal{F}^{-1}[\tilde{\psi}(\mathbf{p})] = \left(\frac{1}{2\pi\hbar}\right)^{3/2} \int_{-\infty}^{\infty} d^3\mathbf{p} e^{i\mathbf{r}\cdot\mathbf{p}/\hbar} \tilde{\psi}(\mathbf{p})$
$\mathcal{F}[\psi^{(n)}(x)] = \left(\frac{ip}{\hbar}\right)^n \tilde{\psi}(p)$	$\tilde{\psi}^{(n)}(p) = \mathcal{F}\left[\left(-\frac{ix}{\hbar}\right)^n \psi(x)\right]$
$\tilde{\psi}(p - p_0) = \mathcal{F}[e^{ip_0x/\hbar} \psi(x)]$	$e^{-ipx_0/\hbar} \tilde{\psi}(p) = \mathcal{F}[\psi(x - x_0)]$
$\mathcal{F}[\psi(cx)] = \tilde{\psi}(p/c)/ c $	$\int_{-\infty}^{\infty} dx \varphi^*(x) \psi(x) = \int_{-\infty}^{\infty} dp \tilde{\varphi}^*(p) \tilde{\psi}(p)$
$\psi(x)$ real: $[\tilde{\psi}(p)]^* = \tilde{\psi}(-p)$	$\psi(x)$ imaginary: $[\tilde{\psi}(p)]^* = -\tilde{\psi}(-p)$
$\Delta x \Delta p \geq \hbar$	

Commutators

Key points

- When a matrix has a block form, we can compute the eigenvalues in each block submatrix.
- The eigenpairs allows you to diagonalize $A = V\Lambda V^{-1}$ in the eigenbasis, where $V = [\mathbf{u}_1|\mathbf{u}_2|\dots]$, $\Lambda = \text{diag}(\lambda_1, \lambda_2, \dots)$, and $A|\mathbf{u}_i\rangle = \lambda_i|\mathbf{u}_i\rangle$. In the eigenbasis we can do $F(A) = VF(\Lambda)V^{-1}$.
- When A is Hermitian, V is unitary: $V^{-1} = V^{\dagger}$.

$ \begin{aligned} [A, B] &= -[B, A] \\ [A, B]^\dagger &= [B^\dagger, A^\dagger] \\ [AB, CD] &= A[B, C]D + AC[B, D] + [A, C]DB + C[A, D]B \\ [A, [B, C]] + [B, [C, A]] + [C, [A, B]] &= 0 \\ e^A e^B &= e^{A+B} e^{\frac{1}{2}[A, B]} \quad ([A, [A, B]] = [B, [A, B]] = 0) \\ [X, P] &= i\hbar \\ [H, P] &= i\hbar \frac{dV(X)}{dX} \end{aligned} $	$ \begin{aligned} [A + B, C + D] &= [A, C] + [A, D] + [B, C] + [B, D] \\ [F(A), A] &= 0 \\ [A, B] = 0 &\implies [F(A), B] = [A, F(B)] = [F(A), F(B)] = 0 \\ [A, [A, B]] = [B, [A, B]] = 0 &\implies [A, F(B)] = [A, B] \frac{dF(B)}{dB} \\ [A, [A, B]] = [B, [A, B]] = 0 &\implies [F(A), B] = [A, B] \frac{dF(A)}{dA} \\ [H, X] &= -\frac{i\hbar}{m} P \\ \langle \varphi_n [A, H] \varphi_n \rangle &= 0, \quad \forall A \end{aligned} $
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- If the matrix is diagonal, the exponential acts directly onto the elements.
- The evolution operator is $U = e^{-iHt/\hbar}$ and it evolves the state by matrix multiplication $U|\psi\rangle$.
- The eigenequation show you the relation of the eigenvectors that must be considered to construct the eigenvectors of the eigenbasis: $A|u_i\rangle = \lambda|u_j\rangle$. Its matrix representation is λ in the ji position.
- You can reduce the dimension of an operator to its eigensubspace when only acting inside it.
- To know the action of an operator you can stimulate it by applying $|\psi\rangle$ or $\langle\psi|$.
- In the operation $|u_i\rangle\langle u_j|$, the element will be located at ij in the matrix.
- Conservative= H time-independent, Stationary state= $|\psi\rangle$ projects in a single eigenstate of H .
- Constant of motion= A time-independent and $[A, H] = 0$.

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