

PHYS 402 (Applications of Quantum Mechanics) Notes

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Introduction:

This set of notes is based on UBC's PHYS 402 (Applications of Quantum Mechanics) course, taught by Dr. Robert Raussendorf. The course covers portions of chapters 1-5 and 7 of Sakurai's "Modern Quantum Mechanics", as well as some additional topics. If any errors are found in the notes, feel free to email me at ryoweil6@student.ubc.ca.

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1 Fundamental Concepts

1.1 The Beginnings of Quantum Mechanics

Before we dive headfirst into the formalism of quantum mechanics, let us first review the first steps of the field as taken in the early 1900s.

Our first founder is Max Planck; the problem at hand was the problem of the blackbody radiation spectrum. The two pre-existing laws (derived from thermodynamics arguments alone) predicting the BBR intensity as a function of wavelength/frequency were flawed. The first was Wien's law (1896):

$$I_{\text{Wien}}(\lambda, T) \sim \frac{1}{\lambda^5} \exp\left(-\frac{1}{\lambda T}\right) \quad (1.1)$$

which agreed with low wavelength/high frequency data well but failed to accurately describe high wavelength/low frequency emission. The second was Rayleigh-Jeans' law (1900):

$$I_{\text{RJ}}(\lambda, T) \sim \frac{T}{\lambda^4} \quad (1.2)$$

which agreed with high wavelength/low frequency data well but failed to accurately describe low wavelength/high frequency emission¹. In fact, the intensity as predicted by Rayleigh-Jeans' diverges at low λ , leading to the (obviously) erroneous conclusion that the total energy emitted by a black body is infinite; the so-called "ultraviolet catastrophe".

In order to solve this problem, in 1900 Planck proposed a quantum hypothesis; that light carries energy in individual packets, or quanta. In particular, for light of frequency f , each quanta carries energy:

$$E = hf. \quad (1.3)$$

Combining this quantum hypothesis with the Boltzmann suppression of high-energy states (from thermodynamics), Planck's law was then derived to be:

$$I_{\text{Planck}}(\lambda, T) = \frac{2hc^2}{\lambda^5} \frac{1}{\exp\left(\frac{hc}{\lambda k_B T}\right) - 1} \quad (1.4)$$

which agrees with the BBR spectrum data across all frequencies². It should also be noted that the integral over all f of the above radiance law yields is finite, resolving the ultraviolet catastrophe. In the above discussion, we have introduced Planck's constant. It has numerical value³:

$$h = 6.626070040 \times 10^{-34} \text{Js} \quad (1.5)$$

h is quantified as "small". What exactly does small mean in this context? For comparison, 1eV is the kinetic energy of an electron acquired in a voltage drop of a Volt, 0.035eV is the average kinetic energy of an atom at room temperature (from $E_k = \frac{3}{2}k_B T$) and 2.4eV is the energy of a single photon from the middle of the visible spectrum (600THz). The energy of a single photon, which depends on h , is in other words "typical" of microscopic phenomena.

Planck's quantum hypothesis would be confirmed in Einstein's (Nobel-prize winning) 1905 explanation of the photoelectric effect (which you likely covered in detail in a previous course in modern physics); namely that quanta of light transfer energy $E = hf$ to electrons in the metal, kicking them out⁴.

¹It should be noted however that a full-derivation of the Rayleigh-Jeans law did not occur until 1905, at which point Planck had already established the more correct explanation.

²Further, we can observe that Planck's law agrees with Wien's law in the high-frequency limit, and with Rayleigh-Jeans' law in the low-frequency limit.

³which is the set/absolute (rather than measured) value of the Planck constant as per the 2018 redefinition of SI units.

⁴Provided of course that $hf > \Phi$ where Φ is the "work function" of the metal.

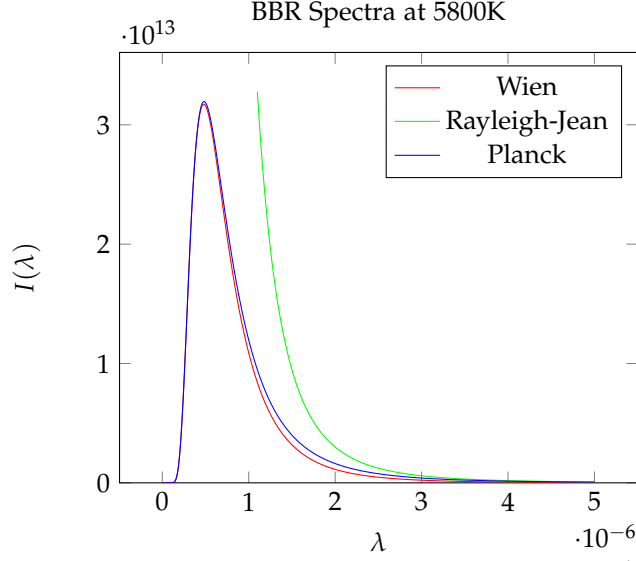


Figure 1.1: Plots of the black body emission spectra at $T = 5800\text{K}$ (the approximate temperature of the surface of the sun) as predicted by Wien’s Law, Rayleigh-Jean’s Law, and Planck’s Law. Planck’s Law was found to agree with experimental observations for all wavelengths. Wien’s Law agrees with observations well in the short wavelength limit but fails for long wavelengths. Rayleigh-Jean’s Law agrees with observations in the long wavelength limit but fails at short wavelengths, and in fact the predicted emitted energy diverges.

Our second founder of interest is DeBroglie. In 1924, he postulated that matter could behave like a wave, positing the DeBroglie wavelength relation:

$$p = \frac{h}{\lambda}. \quad (1.6)$$

The so-called “wave-particle” duality would be confirmed in 1927 by the Davisson-Germer experiment, which saw peaks of electron intensity at distinct angles, showing that electrons scatter in the same nature as photons.

Our third founder of interest is Schrödinger, who postulated the Schrödinger equation (expressed below in the position basis) in 1926:

$$i\hbar \frac{\partial}{\partial t} \psi(\mathbf{r}, t) = \left[\frac{-\hbar^2}{2m} \nabla^2 + V(\mathbf{r}, t) \right] \psi(\mathbf{r}, t). \quad (1.7)$$

It should be noted that this is one of the two core formulas of non-relativistic quantum mechanics, and is the quantum-mechanical equivalent of Newton’s laws. It however does not cover the effects of special relativity (for which we defer the reader to a future course on quantum field theory) or quantum measurement (which we shall address now).

An illuminating demonstration of quantum measurement takes the form of the Stern-Gerlach experiment (first carried out in 1921/1922; see [this article](#) for more historical background). In this experiment, silver atoms are heated and escape from an oven with uniform velocity. The beam of atoms then pass through an inhomogenous magnetic field (generated by an asymmetric pair of magnetic pole pieces) where they are deflected, before hitting a screen where their position is recorded.

Why are silver atoms used for this experiment? Moreover, what exactly is being measured? For this, we consider a simplified model of the atom (which will suffice for the purposes of explaining this experiment).



Figure 1.2: Illustration of the Stern-Gerlach experiment. Silver (Ag) atoms are heated in an oven and escape, and pass through a collimator to form a narrow beam. They then pass through an inhomogeneous magnetic field which deflects the atoms. The position of the atoms is then recorded when they hit the screen.

Silver atoms consist of 47 electrons in the shell, and 47 protons and 61 neutrons in the nucleus. A first guess of the mechanism of the atoms being deflected by the magnetic field may be a Lorentz force effect; however this is not the case as the atoms are electrically neutral. Instead, the silver atom has a single unpaired electron which has an intrinsic angular momentum, known as spin. In particular, the electron is spin-1/2⁵. This provides the silver atom with a net magnetic moment μ proportional to the electron spin⁶ \mathbf{S} :

$$\mu \propto \mathbf{S}. \quad (1.8)$$

We then recall from electromagnetism that a magnetic dipole μ in a magnetic field \mathbf{B} has interaction energy:

$$E = -\mu \cdot \mathbf{B}. \quad (1.9)$$

We can then find the force that the dipole feels by taking the (negative) gradient of the energy:

$$\mathbf{F} = -\nabla(-\mu \cdot \mathbf{B}) = \begin{pmatrix} \frac{\partial}{\partial x} (\mu \cdot \mathbf{B}) \\ \frac{\partial}{\partial y} (\mu \cdot \mathbf{B}) \\ \frac{\partial}{\partial z} (\mu \cdot \mathbf{B}) \end{pmatrix}. \quad (1.10)$$

Ignoring the magnetic fields that are not in the z -direction, we find the force on the silver atoms in the z -direction to be:

$$F_z = \mu_z \frac{\partial B_z}{\partial z}. \quad (1.11)$$

So in the inhomogeneous field produced by the asymmetric magnets, the silver atoms should feel an up/downwards force depending on the direction of \mathbf{S} (which determines μ_z).

Classically, the magnetic moment $\vec{\mu}$ can point in any direction, and therefore μ_z ranges continuously from $+|\mu|$ to $-|\mu|$. Hence, the signature we would expect on the Stern-Gerlach experiment screen (wherein the vertical position of the atoms on the screen corresponds to a measurement of the z -component of the

⁵We will return to a more detailed discussion of angular momentum and spin at a later portion of the course

⁶The astute reader may question why the spin of the unpaired proton in the nucleus has no contribution to the net magnetic moment. This is due to the fact that the proportionality factor between the spin and magnetic moment has a factor of inverse mass. Since the proton is 1836 times heavier than the electron, the proton's magnetic moment contribution is negligible compared to the electron's.



Figure 1.3: The inhomogenous magnetic field used in the Stern-Gerlach experiment, which deflects the silver atoms due to their magnetic dipole moment proportional to electron spin.

magnetic moment) would be a continuous band, as seen in the left of Fig. 1.4 below. However, this is *not* what is observed; instead the experimental result was two discrete dots with nothing in between, as seen in the right of the figure.

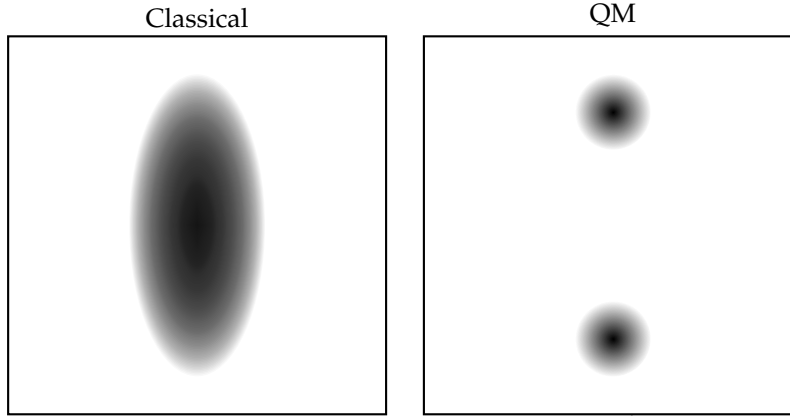


Figure 1.4: Classical prediction (left) and quantum mechanical prediction (right) for the Stern-Gerlach experiment. The screen on the right was observed in experiment.

How do we interpret this result? We can associate the top dot with spins fully polarized upwards (\uparrow) and the bottom dot with spins fully polarized downwards (\downarrow). But why is there no signature for sideways pointing spins? We first will answer how a general spin ($1/2$) state can be represented. If $|\uparrow\rangle$ represents the spin-up state and $|\downarrow\rangle$ represents the spin-down state, then a general spin (and hence sideways spins) can be represented as complex superpositions of these two states, i.e.

$$|\psi\rangle = \alpha|\uparrow\rangle + \beta|\downarrow\rangle \quad (1.12)$$

where $\alpha, \beta \in \mathbb{C}$. What happens in a measurement is then that one element of this general superposition is picked with some probability; indeed, quantum measurement is a probabilistic process. Specifically, we find according to the Born rule that the probability that we measure the spin to be up is $p(\uparrow) = |\alpha|^2$ and the probability that we measure the spin to be down is $p(\downarrow) = |\beta|^2$. Since we require that we measure either spin-up or spin-down, we obtain the normalization condition:

$$p(\uparrow) + p(\downarrow) = |\alpha|^2 + |\beta|^2 = 1. \quad (1.13)$$

The spin state after the measurement is then $|\uparrow\rangle$ or $|\downarrow\rangle$ respectively, according to the Dirac projection

postulate. We will return to these two postulates of quantum mechanics and discuss them in full generality shortly.

However, we will however make a second comment about measurement before concluding this section. Namely, we consider the case where we perform a repeated measurement of the z-component of the spin. As discussed above, the initial general spin state is given by $|\psi\rangle = \alpha|\uparrow\rangle + \beta|\downarrow\rangle$. We then measure the z-component of spin and the post-measurement spin state is $|\uparrow\rangle$ or $|\downarrow\rangle$ with probability $|\alpha|^2$ and $|\beta|^2$ respectively. What happens if we measure the z-component of spin again? We might think that again, we have probability $|\alpha|^2$ of measuring spin-up and probability $|\beta|^2$ of measuring spin-down. But this is *not* the case. If we measured spin-up in the first measurement, we will measure spin-up in the second measurement with probability one. Similarly, if we measured spin-down in the first measurement, we will measure spin-down in the second measurement with probability one. Evidently, the first measurement has done something to the spin such that the measurement probabilities for the second measurement have been affected (they are not the same as the first). This tells us that quantum measurement is a active process that influences the state of the system we measure. Specifically, it is an irreversible process; there is no notion of “undo”-ing the measurement to recover the initial (pre-measurement) state.

1.2 Kets, Bras, and Hilbert Space

Our goal of the initial stages of this course will be to understand the following table:

Quantum states	$ \psi\rangle \in \mathcal{H}$
Evolution	$i\hbar \frac{\partial}{\partial t} \psi\rangle = H \psi\rangle$
Measurement	$ \psi\rangle \mapsto \frac{\Pi_j \psi\rangle}{\sqrt{\langle \psi \Pi_j \psi \rangle}} \quad p(j) = \langle \psi \Pi_j \psi \rangle$

Table 1: Axioms of quantum mechanics, concerning states, evolution, and measurement.

We will discuss the axioms for quantum states and quantum measurement in this chapter, and the axiom for quantum evolution (which readers may recognize as the Schrödinger equation in basis independent form) in the next. It is worth noting that these are the *fundamental postulates* of quantum mechanics; like Newton’s laws of motion in classical mechanics, they cannot be derived. We are only able to interpret them, check if they are consistent, and work out the implications.

Let’s start the axiom for quantum states; after all it will helpful to know what the objects of our interest are, before we start to work with them!

Axiom: Quantum states

Quantum states $|\psi\rangle$ are vectors (also called “kets”) in a complex Hilbert space \mathcal{H} .

The above axiom is only meaningful if we know what a Hilbert space is; its definition is below:

Definition: (Complex) Hilbert spaces

\mathcal{H} is a (complex) Hilbert space if:

- (i) \mathcal{H} is a vector space over \mathbb{C}
- (ii) \mathcal{H} has an inner product
- (iii) \mathcal{H} is complete (with respect to the metric induced by the norm induced by the inner product)⁷-
For the purposes of this course, this last point can be ignored.

Note that the vector space axioms for closure imply that $\forall |\psi\rangle, |\varphi\rangle \in \mathcal{H}$ (where \forall means “for all”) and $\forall c \in \mathbb{C}$, then $|\psi\rangle + |\varphi\rangle \in \mathcal{H}$ and $c|\psi\rangle \in \mathcal{H}$. This tells us that the superposition of quantum states is well defined!

An example which we will return to time and time again (and have already encountered once) is the Hilbert space for a spin-1/2 system. In this case, $\mathcal{H} = \mathbb{C}^2$. A question that may be brooding in the reader’s mind may be “why do we have to use complex numbers?”; one may indeed wonder if real Hilbert spaces may suffice to do quantum mechanics. The response is negative; we indeed need complex numbers! As an illustrative example, consider again the general spin-1/2 state $|\psi\rangle = \alpha|\uparrow\rangle + \beta|\downarrow\rangle$. Suppose we want a state that has equal probability to be measured spin-up or spin-down under a measurement of the z-component of spin. Since $p(\uparrow) = |\alpha|^2$ and $p(\downarrow) = |\beta|^2$, in order to have equal probability we must have $|\alpha| = 1/\sqrt{2}$ and $|\beta| = 1/\sqrt{2}$. A spin pointing in the $+x$ or $-x$ directions indeed has equal weight of up and down. Up to an overall (irrelevant) minus sign, without using complex numbers there are two ways to superimpose $|\uparrow\rangle$ and $|\downarrow\rangle$, from which we can define states corresponding to spins fully polarized in $\pm x$:

$$|x, \pm\rangle = \frac{|\uparrow\rangle \pm |\downarrow\rangle}{\sqrt{2}}. \quad (1.14)$$

However, the $\pm\hat{x}$ and $\pm\hat{y}$ vectors lie in the same $z = 0$ plane, and by symmetry we should require that the $|y, \pm\rangle$ would also have equal weights of $|\uparrow\rangle$ and $|\downarrow\rangle$. But if we limit ourselves to real numbers only, we have already exhausted all possible equal-weight combinations of $|\uparrow\rangle$ and $|\downarrow\rangle$ in Eq. (1.14). We therefore require complex numbers to represent all possible states (and indeed, we find that $|y, \pm\rangle = \frac{|\uparrow\rangle \pm i|\downarrow\rangle}{2}$).

Having motivated the “complex” in the complex vector space part of the definition of Hilbert spaces, let us now motivate the inner product. We want some way to compare quantum states to one another. Our geometric intuition tells us that the states $|\uparrow\rangle$ and $|\nearrow\rangle$ are “close” to each other, while $|\uparrow\rangle$ and $|\downarrow\rangle$ are very “different”. In order to make this intuition rigorous, we define the inner product, and as a prerequisite we define the dual correspondence.

Definition: Dual correspondence

To each vector space \mathcal{H} , there exists a dual vector space \mathcal{H}^* . There is a one-to-one correspondence⁸ between the kets $|\psi\rangle \in \mathcal{H}$ and the bras⁹ $\langle\psi| \in \mathcal{H}^*$. We call this the *dual correspondence*, and write it as follows:

$$|\psi\rangle \overset{DC}{\longleftrightarrow} \langle\psi|. \quad (1.15)$$

It has the following properties:

- (i) $|\psi\rangle + |\varphi\rangle \overset{DC}{\longleftrightarrow} \langle\psi| + \langle\varphi|$
- (ii) $c|\psi\rangle \overset{DC}{\longleftrightarrow} c^* \langle\psi|$

where the $*$ denotes complex conjugation.

Having established the dual correspondence, we may now define the inner product:

⁷This is a technical qualification for the mathematicians in the crowd. An intuitive explanation for the curious; the inner product on a Hilbert spaces creates a notion of distance on the space. There are sequences (of vectors) that get closer together over time; completeness tells us that any such sequences (known as Cauchy sequences) must converge to a limit.

⁹Formally, this follows from the Riesz Representation Theorem. But for the purposes of this course, we take this one-to-one correspondence as a postulate. Curious readers can find discussions/proofs of the theorem in any text on functional analysis, or mathematical quantum theory.

⁹Given such names because $\langle| \rangle$ is a bracket - bra-ket. Physicists remain unmatched in their sense of humour.

Definition: Inner product

We define the inner product between $|\psi\rangle \in \mathcal{H}$ and $|\varphi\rangle \in \mathcal{H}$ as:

$$\langle\varphi|\psi\rangle \in \mathbb{C} \quad (1.16)$$

with the properties:

- (i) $\langle\varphi|\psi\rangle = \langle\psi|\varphi\rangle^*$
- (ii) $\langle\psi|\psi\rangle \geq 0, \forall |\psi\rangle \in \mathcal{H}$
- (iii) $\langle\psi|\psi\rangle = 0 \implies |\psi\rangle = \mathbf{0}$.

Note that $\mathbf{0}$ in the above definition is the null ket (also known as the zero vector, which must be an element of the Hilbert space), where $\mathbf{0} = 0|\psi\rangle, \forall |\psi\rangle$. It is an unphysical state. We normally work with normalized states, i.e. states $|\psi\rangle$ that satisfy $\langle\psi|\psi\rangle = 1$. The null ket has inner product zero and cannot be normalized.

As a first use of the inner product, let us return to our initial motivation for obtaining the “likeness” of states. For normalized states, it follows (and we will later prove) that:

$$0 \leq |\langle\varphi|\psi\rangle| \leq 1. \quad (1.17)$$

We therefore can use the inner product as a method to evaluate the likeness of states. $\langle\varphi|\psi\rangle = 0$ means that $|\psi\rangle$ and $|\varphi\rangle$ are maximally different, and $|\langle\varphi|\psi\rangle| = 1$ corresponds to $|\psi\rangle$ and $|\varphi\rangle$ being the same.

Next, we move onto a discussion of bases of Hilbert spaces. Since Hilbert spaces are vector spaces, they admit a basis. Let us recall what a basis is.

Definition: Basis

A *basis* \mathcal{B} of \mathcal{H} is a set of states $\mathcal{B} = \{|b_j\rangle\}_j$ such that every state $|\psi\rangle \in \mathcal{H}$ can be written in the form:

$$|\psi\rangle = \sum_j \psi_j |b_j\rangle \quad (1.18)$$

with $\psi_j \in \mathbb{C} \forall j$, and the expansion on the RHS is unique.

$|\mathcal{B}|$ is the dimension of \mathcal{H} ¹⁰.

Of particular interest to us will be orthonormal bases, or ONBs.

¹⁰In the case where the Hilbert space is infinite-dimensional, there are additional complications (and in fact, two different kinds of bases!) In general the work we do in this course works perfectly well in the finite-dimensional case, and in the infinite dimensional case we will have to wave our hands a bit in order to avoid getting into the weeds of functional analysis; this is a physics course after all.

Definition: Orthonormal bases

A basis $\mathcal{B} = \{|b_j\rangle\}_j$ is *orthonormal* if

$$\langle b_i | b_j \rangle = \delta_{ij}. \quad (1.19)$$

where δ_{ij} is the Kronecker delta, defined as:

$$\delta_{ij} = \begin{cases} 1 & i = j \\ 0 & i \neq j \end{cases}. \quad (1.20)$$

One can obtain the expansion coefficients ψ_j with respect to ONBs. Writing $|\psi\rangle = \sum_j \psi_j |b_j\rangle$ as in Eq. (1.18) and taking the inner product of $|\psi\rangle$ with a vector $|b_i\rangle$ from the ONB, we have:

$$\langle b_i | \psi \rangle = \sum_j \psi_j \langle b_i | b_j \rangle = \sum_j \psi_j \delta_{ij} = \psi_i. \quad (1.21)$$

We will now prove a useful trick involving ONBs.

Theorem: Resolution of the identity

For all ONBs $\{|b_j\rangle\}_j$, the following relation holds:

$$\sum_j |b_j\rangle \langle b_j| = \mathbb{I} \quad (1.22)$$

where \mathbb{I} is the identity operator on the Hilbert space.

Proof. Recall that $|\psi\rangle = \sum_j \psi_j |b_j\rangle$ for any $|\psi\rangle \in \mathcal{H}$ and for any basis $\{|b_j\rangle\}_j$ of \mathcal{H} . Further, recall that $\psi_j = \langle b_j | \psi \rangle$ if the basis is orthonormal. Hence we have that:

$$|\psi\rangle = \sum_j \langle b_j | \psi \rangle |b_j\rangle = \sum_j |b_j\rangle \left(\langle b_j | \psi \rangle \right) = \sum_j \left(|b_j\rangle \langle b_j| \right) |\psi\rangle = \left(\sum_j |b_j\rangle \langle b_j| \right) |\psi\rangle. \quad (1.23)$$

Since the above relation holds for all $|\psi\rangle$, it follows then that $\sum_j |b_j\rangle \langle b_j|$ is the identity as claimed. \square

At this point in the course, the reader may be wondering what happened to quantum *wavefunctions*¹¹; the central objects of interest have instead been quantum states, without a wavefunction in sight. We now elucidate the connection between the two.

¹¹Though this nomenclature of “wavefunction” is arguably a misnomer; the Schrödinger equation does not contain second order derivatives in time, as a wave equation would.

Definition: Wavefunctions

Consider (as before) expanding $|\psi\rangle$ in a ONB $\{|b_j\rangle\}_j$. We then have:

$$|\psi\rangle = \sum_j |b_j\rangle \langle b_j|\psi\rangle. \quad (1.24)$$

Therein, $\psi_j = \psi(j) = \langle b_j|\psi\rangle$ is the *wavefunction*; a wavefunction is just a quantum state expanded in a given ONB.

A particularly important (and familiar) example is using the position basis. Consider the resolution of the identity involving position:

$$\int dx |x\rangle \langle x| = \mathbb{I}. \quad (1.25)$$

For any $|\psi\rangle$, we then have:

$$|\psi\rangle = \int dx |x\rangle \langle x|\psi\rangle. \quad (1.26)$$

Where $\langle x|\psi\rangle = \psi(x)$ is the position wavefunction, which played a central role in your first course in quantum mechanics. However, the reader should now recognize that $|\psi\rangle$ is a more fundamental object than this wavefunction, as it not only contains the information for $\psi(x)$ but also for $\tilde{\psi}(p)$ (the momentum wavefunction) or any other wavefunction; any wavefunction is just the expansion coefficients of the state in a given basis.

We have now established quite a bit of machinery to discuss quantum states, but have not done anything with the states; let's change that by discussing measurements! In a quantum measurement, the alternatives $\{|m_j\rangle, j \in \text{Outcomes}\}$ for the states after measurement will:

- Form a basis
- And are pairwise orthogonal, i.e. $\langle m_i|m_j\rangle = 0, \forall i \neq j$.

With the necessary mathematical formalism under our belt, we can now state a first version of the Dirac postulate and Born rule axioms.

Axiom: Quantum measurement (version 1)

Dirac postulate: Under quantum measurement, the measured quantum state $|\psi\rangle$ is probabilistically changed into one of a number of alternatives $\{|m_j\rangle\}_j$, with:

$$\langle m_i|m_j\rangle = \delta_{ij}. \quad (1.27)$$

Note that $\{|m_j\rangle\}_j$ forms an ONB.

Born rule: Given a quantum state $|\psi\rangle$, the probability for obtaining outcome j , corresponding to post-measurement state $|m_j\rangle$, is:

$$p_j = |\langle m_j|\psi\rangle|^2. \quad (1.28)$$

There are now two questions that may arise. The first is that the statement of the Dirac projection postulate and the Born rule do not match that found in Table 1. The second question concerns the post-measurement states $\{|m_j\rangle\}_j$; namely, how are they determined? We will address the latter question first, and build up the formalism to state the measurement axioms in full generality. To this end, we move to a discussion of linear operators, which describe quantum mechanical observables and evolution.

1.3 Operators and Observables

Definition: Linear Operators

A is an *operator* that acts on a Hilbert space \mathcal{H} if $\forall |\psi\rangle \in \mathcal{H}, A|\psi\rangle \in \mathcal{H}$. A is *linear* if:

$$A(\alpha|\psi\rangle + \beta|\varphi\rangle) = \alpha(A|\psi\rangle) + \beta(A|\varphi\rangle) \quad (1.29)$$

$\forall |\psi\rangle, |\varphi\rangle \in \mathcal{H}$ and $\forall \alpha, \beta \in \mathbb{C}$.

A point of notation; we will use capital letters to denote operators in this course. Some sources also use hats to denote this (e.g. \hat{A}). Note that linear operators can be added and multiplied (more accurately, composed) to yield other linear operators. They are associative and distributive under these operations, i.e. for all linear operators A, B, C we have:

$$(A + B) + C = A + (B + C) \quad (1.30)$$

$$(AB)C = A(BC) \quad (1.31)$$

$$A(B + C) = AB + AC. \quad (1.32)$$

However, note that in general operators are *not* commutative, that is:

$$AB \neq BA. \quad (1.33)$$

Quantum mechanical observables are a specific type of linear operators; namely, they are Hermitian. In order to make sense of this, we introduce two more definitions.

Definition: Hermitian adjoint

The *Hermitian adjoint* A^\dagger of a linear operator A is defined via the dual correspondence:

$$A|\psi\rangle \xleftrightarrow{DC} \langle\psi|A^\dagger. \quad (1.34)$$

Definition: Hermitian operators and observables

An operator A is *Hermitian* if:

$$A^\dagger = A. \quad (1.35)$$

Quantum mechanical *observables* (such as position, momentum, spin, and energy) are Hermitian operators.

In general, acting on a state with an operator changes the state, and the new state is not necessarily proportional to the original state, i.e:

$$A|\psi\rangle \not\propto |\psi\rangle. \quad (1.36)$$

However, for special states known as eigenstates of an operator, this is indeed the case.

Definition: Eigenstates and eigenvalues

An *eigenstate* $|a\rangle$ of a linear operator A is a state that satisfies:

$$A|a\rangle = a|a\rangle. \quad (1.37)$$

Therein, $a \in \mathbb{C}$ is the *eigenvalue* corresponding to that eigenstate. The null vector $\mathbf{0}$ is excluded from being an eigenvector.

Having defined these objects, we state (but do not prove) an important theorem concerning observables:

Theorem: Diagonalization/Spectral Theorem

A Hermitian operator A (i.e. any observable) can be diagonalized; that is, it is able to be written as:

$$A = \sum_i a_i |a_i\rangle \langle a_i| \quad (1.38)$$

Where the eigenvectors $\{|a_i\rangle\}$ of A forms an orthonormal basis¹²(an *eigenbasis*) and a_i are the corresponding eigenvalues.

With some more machinery developed, we can more carefully state the Dirac postulate and the Born rule.

Axiom: Quantum measurement (version 2)

Dirac postulate:

1. Each possible outcome of the measurement of an observable A is an eigenvalue of A .
2. If the eigenvalue a is found in the measurement, then the post measurement state is an eigenvector $|a\rangle$ of A ,

$$|\psi\rangle \mapsto |a\rangle \quad (1.39)$$

satisfying

$$A|a\rangle = a|a\rangle. \quad (1.40)$$

Born rule: Given an initial state of $|\psi\rangle$, the possibility for the measurement outcome a_i occurring in a measurement of an observable A (where a_i is an eigenvalue of A) is:

$$p_i = |\langle a_i | \psi \rangle|^2. \quad (1.41)$$

We reiterate that the eigenstate $|a\rangle$ is the possible post-measurement state, and the eigenvalue a is the corresponding measurement outcome. As a concrete example, the spin- z operator S_z (for spin-1/2) systems has eigenstates $|\uparrow\rangle$ and $|\downarrow\rangle$, with corresponding eigenvalues of $+\hbar/2$ and $-\hbar/2$ (i.e. $S_z|\uparrow\rangle = +\hbar/2|\uparrow\rangle$, and likewise for $|\downarrow\rangle$). We may write S_z in diagonal form as:

$$S_z = \frac{\hbar}{2}(|\uparrow\rangle\langle\uparrow| - |\downarrow\rangle\langle\downarrow|). \quad (1.42)$$

A measurement of the z component of spin has two possible outcomes $\pm\hbar/2$, corresponding to post measurement states $|\uparrow\rangle/|\downarrow\rangle$. $\{|\uparrow\rangle, |\downarrow\rangle\}$ is an ONB, and we can expand a general state $|\psi\rangle$ in this basis as $|\psi\rangle = \alpha|\uparrow\rangle + \beta|\downarrow\rangle$. The Born rule then tells us that $p(\uparrow) = |\langle\uparrow|\psi\rangle|^2 = |\alpha|^2$ and similarly that $p(\downarrow) = |\beta|^2$.

There are two points of consistency that the restatement of the Dirac postulate invites. First, an experiment should only have real-valued outcomes; an measurement shouldn't return a complex number. Second, it is not a priori immediate that the eigenstates/post-measurement states are mutually orthogonal, as the first statement of the postulate requires. Fortunately, there is a theorem that covers both.

¹²If the eigenvalues of A are non-degenerate, then the set of eigenstates of A are automatically mutually orthogonal, as we prove in the next theorem. If A has degenerate eigenvalues, two eigenstates with the same eigenvalue are not necessarily guaranteed to be orthogonal, but we are still able to choose eigenstates (picking orthogonal vectors from the degenerate subspaces) to form an ONB out of the eigenstates; so we need not worry too much.

Theorem

The eigenvalues of a Hermitian operator A are real, and the eigenvectors of A corresponding to distinct eigenvalues are orthogonal.

Proof. Consider two eigenstates $|a\rangle, |a'\rangle$ of A (with corresponding eigenvalues a, a'). First, we make the observation that $\langle a|A = \langle a|a^*$. This follows as:

$$\langle a|a^* \xleftrightarrow{DC} a|a\rangle = A|a\rangle = A^\dagger|a\rangle \xleftrightarrow{DC} \langle a|A \quad (1.43)$$

where the first equality invokes that $|a\rangle$ is an eigenvector of A , and the second equality invokes the Hermiticity of A . Since the dual correspondence is unique, we can compare the first and the last expressions to conclude that $\langle a|A = \langle a|a^*$. Next, we consider the number $\langle a|A|a'\rangle$. We may write this as:

$$\langle a|A|a'\rangle = \langle a|(A|a'\rangle) = \langle a|a'|a\rangle = a'\langle a|a'\rangle \quad (1.44)$$

but by associativity, we can just as well write this as:

$$\langle a|A|a'\rangle = (\langle a|A)|a'\rangle = \langle a|a^*|a'\rangle = a^*\langle a|a'\rangle. \quad (1.45)$$

Therefore we find that $a^*\langle a|a'\rangle = a'\langle a|a'\rangle$. From this we obtain that:

$$(a^* - a')\langle a|a'\rangle = 0. \quad (1.46)$$

There are now two cases to consider.

- (I) If $|a\rangle = |a'\rangle$, then $\langle a|a'\rangle = \langle a|a\rangle > 0$ (as $|a\rangle$ is an eigenvector, it cannot be a null vector). Therefore for Eq. (1.46) to be satisfied it must follow that $a^* = a$, i.e. a is real.
- (II) If instead $a \neq a'$, then $a^* \neq a'$ (as $a = a^*$), so for Eq. (1.46) to be satisfied it must follow that $\langle a|a'\rangle = 0$.

□

We have now shown that the Dirac projection postulate is consistent with what we should expect out of measurements. However, we now clarify what is often a point of confusion, that being the difference between individual and averaged measurement outcomes. The dirac postulate states that given a state $|\psi\rangle$, a possible outcome of the measurement of an observable A are eigenvalues a of A . This speaks to possible outcomes in *individual measurements*. The *expectation value* is conceptualized quite differently.

Definition: Expectation values

The *expectation value* $\langle A \rangle_\psi$ is the average outcome in the measurement of A on $|\psi\rangle$, i.e.:

$$\langle A \rangle_\psi := \sum_i p_i a_i \quad (1.47)$$

where p_i is the probability of measuring outcome a_i . We are able to write Eq. (1.47) equivalently as:

$$\langle A \rangle_\psi = \sum_i |\langle a_i|\psi\rangle|^2 a_i = \sum_i \langle a_i|\psi\rangle^* \langle a_i|\psi\rangle a_i = \sum_i \langle \psi|a_i\rangle a_i \langle a_i|\psi\rangle = \langle \psi| \left(\sum_i a_i |a_i\rangle \langle a_i| \right) |\psi\rangle = \langle \psi|A|\psi\rangle. \quad (1.48)$$

where in the first equality we use the Born rule, and in the last equality we consider A in diagonalized form.

The formalism of quantum mechanics is therefore well-suited to predict both individual and averaged measurement outcomes; but take care not to confuse them! In general $\langle A \rangle_\psi$ is *not* a possible (individual) measurement outcome.

Up until now, we have considered both kets $|\psi\rangle$ and operators A as abstract objects; however sometimes it is helpful to cast these into a more concrete form in order to do computations. We thus introduce the idea of matrix/vector representations.

Consider an ONB $\mathcal{B} = \{|b_i\rangle\}_i$, and consider an expression of the form $|\varphi\rangle = A|\psi\rangle$. Inserting the resolution of the identity once on the left hand side and twice on the right hand side, we have:

$$\left(\sum_i |b_i\rangle\langle b_i|\right) |\varphi\rangle = \left(\sum_i |b_i\rangle\langle b_i|\right) A \left(\sum_j |b_j\rangle\langle b_j|\right) |\psi\rangle \quad (1.49)$$

Now redrawing some brackets we obtain:

$$\sum_i |b_i\rangle (\langle b_i|\varphi\rangle) = \sum_i \sum_j |b_i\rangle (\langle b_i|A|b_j\rangle) (\langle b_j|\psi\rangle). \quad (1.50)$$

Now, we can consider $\langle b_i|\varphi\rangle = [\varphi]_i$ and $\langle b_j|\psi\rangle = [\psi]_j$ (where we use $[\cdot]$ to denote ‘representation of’) as elements of column vectors $[\varphi]$ and $[\psi]$, and $\langle b_i|A|b_j\rangle = [A]_{ij}$ as elements of a matrix $[A]$. In other words, expanded out in the $\{|b_i\rangle\}_i$ basis, we can realize $[\varphi]$ is the column vector obtained by multiplying the column vector $[\psi]$ by the matrix $[A]$.

Similarly, we can write for any bra $\langle\tau|$ that:

$$\langle\tau| = \langle\tau|\sum_i |b_i\rangle\langle b_i| = \sum_i \langle\tau|b_i\rangle\langle b_i| = \sum_i [\tau]_i \langle b_i|. \quad (1.51)$$

This yields us the following way of thinking about bras/kets/operators:

Abstract Objects	Representation in ONB
Ket $ \varphi\rangle$	Column vector $[\varphi]$
Operator A	Matrix $[A]$
Bra $\langle\tau $	Row vector $[\tau]$

Table 2: Abstract objects and their representations when expanded out in a basis.

However, do take note that a ket is *not* equal to a column vector. The column vector is a representation of a ket, much in the same way that $\begin{pmatrix} 1 \\ 0 \end{pmatrix}$ is not in itself a vector but instead a concrete (standard) representation of the abstract vector e_1 .

Let us supplement this discussion of matrix representations by returning to our favourite example of spin-1/2. Consider the ONB $\mathcal{B} = \{|\uparrow\rangle, |\downarrow\rangle\}$ of $\mathcal{H} = \mathbb{C}^2$. Making the identification that $|\uparrow\rangle \cong \begin{pmatrix} 1 \\ 0 \end{pmatrix}$ and $|\downarrow\rangle \cong \begin{pmatrix} 0 \\ 1 \end{pmatrix}$, for any $|\psi\rangle \in \mathcal{H}$ we then have that:

$$|\psi\rangle \cong \begin{pmatrix} \langle\uparrow|\psi\rangle \\ \langle\downarrow|\psi\rangle \end{pmatrix} = [\psi] \quad (1.52)$$

and for any linear operator A acting on kets in \mathcal{H} we have:

$$A \cong \begin{pmatrix} \langle\uparrow|A|\uparrow\rangle & \langle\uparrow|A|\downarrow\rangle \\ \langle\downarrow|A|\uparrow\rangle & \langle\downarrow|A|\downarrow\rangle \end{pmatrix} = [A]. \quad (1.53)$$

Taking spin-z operator S_z as a concrete example, we can represent it in this basis as:

$$S_z = \frac{\hbar}{2} (|\uparrow\rangle\langle\uparrow| - |\downarrow\rangle\langle\downarrow|) \cong \frac{\hbar}{2} \begin{pmatrix} 1 & 0 \\ 0 & -1 \end{pmatrix}. \quad (1.54)$$

As another example, consider the spin- x operator S_x . It has eigenstates:

$$|\pm\rangle := \frac{|\uparrow\rangle \pm |\downarrow\rangle}{\sqrt{2}} \quad (1.55)$$

we can represent it as:

$$S_x = \frac{\hbar}{2} (|+\rangle\langle+| - |-\rangle\langle-|) \cong \frac{\hbar}{2} \begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}. \quad (1.56)$$

From the work we have done so far, it may have become clear that it is generally easiest to work in the eigenbasis of whatever observable is being considered. For example, if considering the measurement of the observable S_z , it is natural to work with the basis $\mathcal{B} = \{|\uparrow\rangle, |\downarrow\rangle\}$. This fact will continue to be true when we later consider Schrödinger evolution of observables. However, there are times when there are multiple observables under consideration; in the measurement picture, this could be the sequential measurement of observable O_1 followed by the measurement of observable O_2 ¹³. In this case, we have two different bases of interest, namely the eigenbasis of O_1 and the eigenbasis of O_2 . We may then find it useful to consider a transformation between these bases; let us work through that now.

Consider two ONBs defined by $\mathcal{B} = \{|b_i\rangle\}_i$ and $\mathcal{B}' = \{|a_j\rangle\}_j$, and some state $|\psi\rangle \in \mathcal{H}$. Inserting the resolution of identity $\mathbb{I} = \sum_j |a_j\rangle\langle a_j|$ between the wavefunction $\langle b_i|\psi\rangle = \psi_i$, we obtain:

$$\langle b_i|\psi\rangle = \langle b_i| \left(\sum_j |a_j\rangle\langle a_j| \right) |\psi\rangle = \sum_j \langle b_i|a_j\rangle \langle a_j|\psi\rangle \quad (1.57)$$

Now, $\langle b_i|\psi\rangle$ can be viewed as elements of a column vector $[\psi]$ (whose components are the expansion coefficients of $|\psi\rangle$ in the basis \mathcal{B}), $\langle a_i|\psi\rangle$ can be analogously viewed as the elements of a column vector $[\tilde{\psi}]$ (whose components are the expansion coefficients of $|\psi\rangle$ in the basis \mathcal{B}'), and $\langle b_i|a_j\rangle = T_{ij}$ can be viewed as the matrix elements of a (matrix) representation of the transformation operator T between the two bases. We can therefore write Eq. (1.57) as a matrix/vector equation:

$$[\psi] = [T][\tilde{\psi}]. \quad (1.58)$$

The operator T actually has a special property; namely, it is *unitary*.

Definition: Unitary operators

A linear operator U is *unitary* if:

$$UU^\dagger = U^\dagger U = \mathbb{I}. \quad (1.59)$$

Theorem

The basis transformation operator between bases $\mathcal{B} = \{|b_i\rangle\}_i$ and $\mathcal{B}' = \{|a_j\rangle\}_j$ defined as:

$$T = \sum_n |a_n\rangle\langle b_n| \quad (1.60)$$

which has matrix representation (in either \mathcal{B} or \mathcal{B}'):

$$[T]_{ij} = \langle b_i|a_j\rangle \quad (1.61)$$

is unitary.

¹³An similar setting of interest with Schrödinger evolution could be the evolution of a quantum state under a Hamiltonian H_1 for some time $[t_0, t_1]$ followed by evolution by a different Hamiltonian H_2 for some time $[t_1, t_2]$.

Proof. First, it should be clear from the basis agnostic definition in Eq. (1.60) that T is in fact a basis transformation operator, as by inspection it satisfies $T|b_n\rangle = |a_n\rangle$ for $n = 1, 2, \dots, d$. We next verify that it indeed has the claimed matrix representation; in the \mathcal{B} basis, we have:

$$[T]_{ij} = \langle b_i | T | b_j \rangle = \langle b_i | \left(\sum_n |a_n\rangle \langle b_n| \right) | b_j \rangle = \sum_n \langle b_i | a_n \rangle \langle b_n | b_j \rangle = \sum_n \langle b_i | a_n \rangle \delta_{nj} = \langle b_i | a_j \rangle \quad (1.62)$$

and in the \mathcal{B}' basis we have:

$$[T]_{ij} = \langle a_i | T | a_j \rangle = \langle a_i | \left(\sum_n |a_n\rangle \langle b_n| \right) | a_j \rangle = \sum_n \langle a_i | a_n \rangle \langle b_n | a_j \rangle = \sum_n \delta_{in} \langle b_n | a_j \rangle = \langle b_i | a_j \rangle \quad (1.63)$$

so Eq. (1.61) indeed holds as claimed.

We move onto the unitarity proof. We claim that the equality below:

$$[T^\dagger]_{kl} = [T]_{lk}^* \quad (1.64)$$

holds *in general* for any linear operator T . Note that Eq. (1.64) reconciles the abstract notion of the Hermitian adjoint with the familiar operation of taking the conjugation transpose of a matrix (which you should have encountered in your linear algebra course). The equation follows from the definition of the Hermitian adjoint using the dual correspondence, and is left as an exercise for the reader.

Applying Eq. (1.64) to the definition of T , we find:

$$[T^\dagger]_{kl} = [T]_{lk}^* = \langle b_l | a_k \rangle^* = \langle a_k | b_l \rangle. \quad (1.65)$$

Next, we consider the matrix elements of the operator $T^\dagger T$:

$$[T^\dagger T]_{kj} = \sum_l [T^\dagger]_{kl} [T]_{lj} = \sum_l \langle a_k | b_l \rangle \langle b_l | a_j \rangle = \langle a_k | \left(\sum_l |b_l\rangle \langle b_l| \right) | a_j \rangle. \quad (1.66)$$

where in the first equality we use the definition of matrix multiplication, and in the second equality we use Eq. (1.65). Recognizing the resolution of the identity, the above equation then becomes:

$$[T^\dagger T]_{kj} = \langle a_k | \mathbb{I} | a_j \rangle = \langle a_k | a_j \rangle = \delta_{kj} \quad (1.67)$$

where in the last equality we use that $\{|a_j\rangle\}_j$ is an ONB. From $[T^\dagger T]_{kj} = \delta_{kj}$ we can conclude that $T^\dagger T = \mathbb{I}$, and $TT^\dagger = \mathbb{I}$ can be shown analogously. We conclude that T is unitary. \square

Note that while this is our first encounter with unitarity, it will certainly not be our last; unitary operators will have a very large role to play when we consider Schrödinger evolution.

As an example, let us solve for T which transforms from the S_z eigenbasis $\mathcal{B} = \{|\uparrow\rangle, |\downarrow\rangle\}$ to the S_x eigenbasis $\mathcal{B}' = \{|+\rangle = \frac{|\uparrow\rangle + |\downarrow\rangle}{\sqrt{2}}, |-\rangle = \frac{|\uparrow\rangle - |\downarrow\rangle}{\sqrt{2}}\}$. We may write it down in operator form as:

$$T = |+\rangle \langle \uparrow| + |-\rangle \langle \downarrow|. \quad (1.68)$$

To get a clearer picture, we can write down its matrix representation. Computing the inner products between the eigenvectors of S_z and S_x , we obtain:

$$[T] = \begin{pmatrix} \langle \uparrow | + \rangle & \langle \uparrow | - \rangle \\ \langle \downarrow | + \rangle & \langle \downarrow | - \rangle \end{pmatrix} = \begin{pmatrix} \frac{1}{\sqrt{2}} & \frac{1}{\sqrt{2}} \\ \frac{1}{\sqrt{2}} & -\frac{1}{\sqrt{2}} \end{pmatrix} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}. \quad (1.69)$$

The above matrix representation is the same whether we choose the identification $|\uparrow\rangle \cong \begin{pmatrix} 1 \\ 0 \end{pmatrix}, |\downarrow\rangle \cong \begin{pmatrix} 0 \\ 1 \end{pmatrix}$ or $|+\rangle \cong \begin{pmatrix} 1 \\ 0 \end{pmatrix}, |-\rangle \cong \begin{pmatrix} 0 \\ 1 \end{pmatrix}$ (but note: this is NOT true of operators in general. For example the matrix representations of S_z and S_x will look different depending on the identifications chosen). It can be easily verified via matrix multiplication that T is unitary, which is consistent with the general theorem we just proved. The T above is a ubiquitous operator in the field of quantum computation, and is called the *Hadamard operator*.

1.4 Projectors and Measurement

We now return to the setting of measurement to resolve an ambiguity that was left by version 2 of our measurement axioms. Namely, we ask “what if the observable in question has degenerate eigenvalues?” (Degenerate referring to the fact that two or more distinct eigenstates of an observable may have the same eigenvalues). To see why this is a problem with our current formulation, consider an operator A with two distinct eigenstates $|a\rangle, |\tilde{a}\rangle$ such that:

$$A|a\rangle = a|a\rangle, A|\tilde{a}\rangle = a|\tilde{a}\rangle. \quad (1.70)$$

In our earlier formulation of the Dirac postulate, we claimed that a measurement of an observable A with outcome/eigenvalue a would change the quantum state to be an eigenstate with the corresponding eigenvalue. Now that there are two distinct eigenstates with the same eigenvalue, which eigenstate is chosen?

As a concrete example, we consider the z -component of spin for a spin-1 particle. The S_z operator has three eigenstates, namely $|+\rangle, |0\rangle$ (this is *not* the null ket/zero vector!) and $|-\rangle$ which forms a basis, and can be written as:

$$S_z = \hbar(|+\rangle\langle+| - |-\rangle\langle-|). \quad (1.71)$$

$|+\rangle$ has eigenvalue $+\hbar$, $|0\rangle$ has eigenvalue zero, and $|-\rangle$ has eigenvalue $-\hbar$. Now, consider the observable S_z^2 . From S_z above, we can deduce this to be:

$$S_z^2 = \hbar^2(|+\rangle\langle+| + |-\rangle\langle-|) \quad (1.72)$$

Now, we find that both $|+\rangle$ and $|-\rangle$ have eigenvalue $+\hbar^2$. So, if we measured S_z^2 for a state $|\psi\rangle$ and got outcome \hbar^2 , our current formulation is ill-equipped to deduce what the post-measurement state would be. In order to refine our formulation, we require the notion of a projector.

Another example (which we cannot discuss in full detail yet) is that of “spooky action at a distance” (coined by the one and only Einstein). In this setting, we have a two-particle system which is in an entangled Bell state (to be defined in full mathematical detail later; for now, take entangled to mean “stronger than classically correlated”, or “it is impossible to describe the individual subsystems/particles, and only the joint system can be described”). We then separate the two particles some great distance; say the Earth and moon. When we measure the particle on Earth (but do nothing to the particle on the moon), we find that upon knowing the Earth measurement outcome we *immediately* know what the post-measurement state of the particle on the moon is. This seems counterintuitive, and indeed, superluminal; it seems as though the measurement on Earth affects the state on the moon faster than the speed of light! While this setup does not actually turn out to violate special relativity (and indeed, we will prove later in the course that no quantum communication protocol can be superluminal), it is certainly of interest as depicting non-classical phenomenology. It turns out that the measurement of just the particle on Earth conducted in this setting corresponds to a measurement of a degenerate (joint) observable; we therefore also require a reformulation of our measurement axioms involving projectors here. Without further ado, let us define what a projector is.



Figure 1.5: Visualization of the “spooky action at a distance” scenario. A two-particle state is prepared in an entangled Bell state. The particles are then spatially separated a great distance; here depicted as one particle on the Earth and one particle in the moon. When the particle on Earth (and *only* the particle on Earth) is measured, the entanglement structure of the state leads to the particle on the moon being projected into a post-measurement state that is immediately known to the experimenter on Earth based on their measurement outcome. The measurement conducted in this scenario of interest turns out to be one with degeneracy, and hence we require a reformulation of measurement in order to describe it.

Definition: Projectors

A linear operator Π is a *projector* if it satisfies:

$$\Pi^2 = \Pi^\dagger = \Pi. \quad (1.73)$$

Below are examples of projectors in matrix representations:

$$\Pi_1 \cong \begin{pmatrix} 1 & 0 & 0 \\ 0 & 0 & 0 \\ 0 & 0 & 0 \end{pmatrix}, \quad \Pi_2 \cong \begin{pmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 0 \end{pmatrix}. \quad (1.74)$$

Π_1 is a rank 1 projector, while Π_2 is rank 2 > 1 . Why we call an operator with the properties in Eq. (1.73) a projector might not be obvious, but the nomenclature is elucidated by the above examples. A projector projects a state into a lower-dimensional subspace of the Hilbert space. Π_1 has the property of taking a three-dimensional vector and projecting it into a 1-dimensional subspace, while Π_2 has the property of taking a three-dimensional vector and projecting it into a 2-dimensional subspace. Π is a projector (though a trivial one), and is a projection from a space to itself. In Fig. 1.6 we visualize the action of Π_1, Π_2 for the case when our vector space is \mathbb{R}^3 (but one should keep in mind that this is for the sake of intuition, and the Hilbert spaces we use in quantum mechanics are, of course, complex).

Now let's return to the bra-ket formalism and what projectors look like in this abstract setting. First, recall that we can write an observable A in the form:

$$A = \sum_i a_i |a_i\rangle \langle a_i| \quad (1.75)$$

where $|a_i\rangle$ is the eigenstate of A corresponding to the eigenvalue a_i , and $\{|a_i\rangle\}_i$ is an ONB. In the non-degenerate case, each of the a_i s are distinct. However, in general degenerate eigenvalues (where $a_i = a_j$ for some i, j) are possible, and the current form of the expression does not make this particularly clear. With our knowledge of projectors, let us now rewrite the above as:

$$A = \sum_a a \Pi_a \quad (1.76)$$

where each of the a s are *distinct* eigenvalues of A , and

$$\Pi_a = \sum_{a_i=a} |a_i\rangle \langle a_i| \quad (1.77)$$

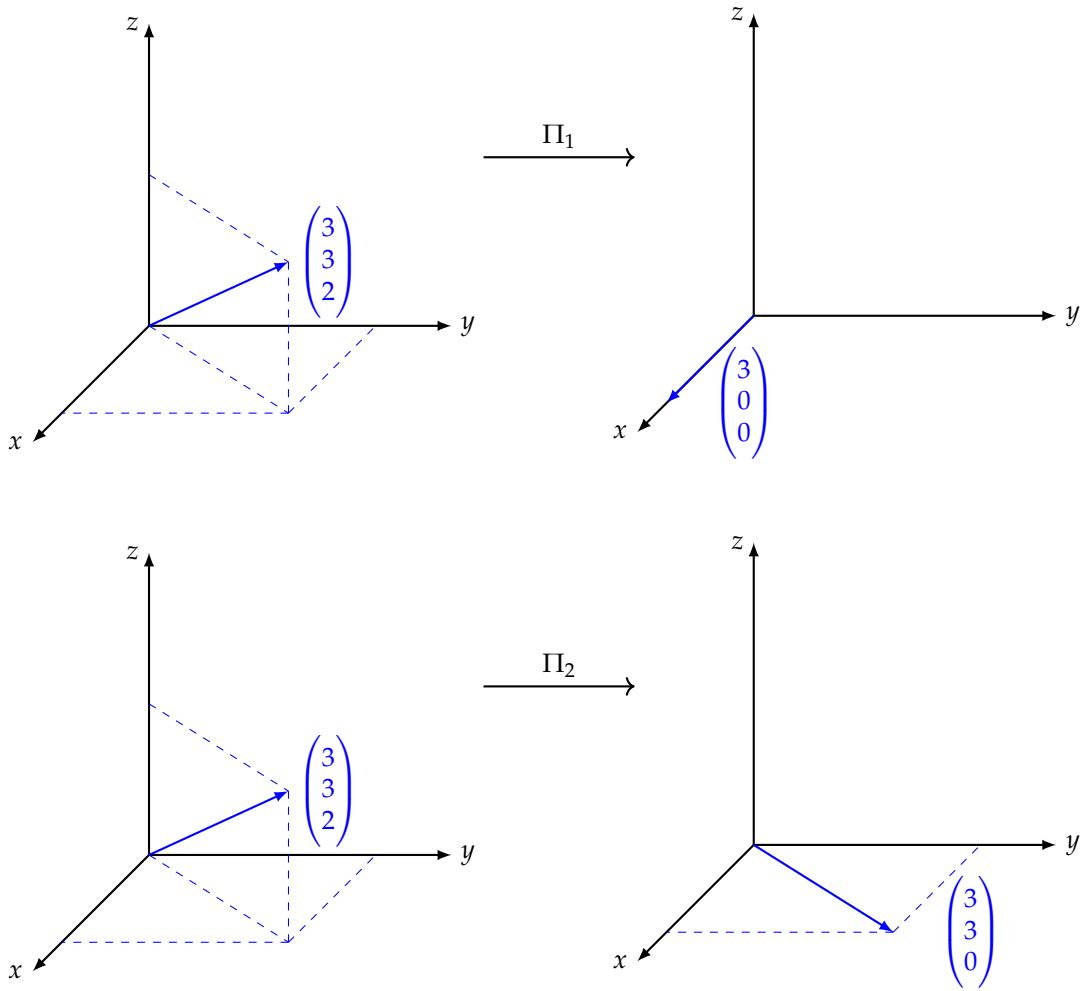


Figure 1.6: Visualization of the action of projectors Π_1, Π_2 (as defined in Eq. (1.74)) on a vector in \mathbb{R}^3 . Π_1 can be visualized as projecting the given vector onto the one-dimensional subspace that is the x -axis subspace; preserving the x -component of the vector, and nullifying the y and z -components. Π_2 can be visualized as projecting the given vector onto the two-dimensional subspace that is the xy -plane; preserving the x and y components of the vector and nullifying the z -component.

is the *projector onto the eigenvalue- a subspace*. Let us verify that these are indeed projectors. First, we verify that they are Hermitian:

$$\Pi_a^\dagger = \left(\sum_{a_i=a} |a_i\rangle\langle a_i| \right)^\dagger = \sum_{a_i=a} (|a_i\rangle\langle a_i|)^\dagger = \sum_{a_i=a} |a_i\rangle\langle a_i| = \Pi_a. \quad (1.78)$$

where in the second-to-last equality we use that $(|a\rangle\langle b|)^\dagger = |b\rangle\langle a|$ (which follows immediately from the definition of the Hermitian adjoint; the proof is left to the reader!). Next, we show that they are idempotent (that is, they square to themselves):

$$\Pi_a^2 = \left(\sum_{a_i=a} |a_i\rangle\langle a_i| \right)^2 = \sum_{a_i=a} \sum_{a_j=a} |a_i\rangle\langle a_i|a_j\rangle\langle a_j| = \sum_{a_i=a} \sum_{a_j=a} |a_i\rangle\langle a_j|\delta_{ij} = \sum_{a_i=a} |a_i\rangle\langle a_i| = \Pi_a. \quad (1.79)$$

So they are indeed projectors! In this form, we have decomposed the observable A into the parts associated with each eigenvalue in a clear way. These projectors have some properties of note, described in the theorem below.

Theorem

Let $\{\Pi_a\}_a$ be the set of projectors associated to an observable A (with $\Pi_a = \sum_{a_i=a} |a_i\rangle\langle a_i|$ being the projector onto the eigenvalue- a subspace). These projectors are mutually orthogonal:

$$\Pi_i\Pi_j = \delta_{ij}\Pi_i \quad (1.80)$$

and are complete:

$$\sum_a \Pi_a = \mathbb{I}. \quad (1.81)$$

Proof. The idempotency of projectors covers the $i = j$ case in Eq. (1.80), and if $i \neq j$, then the expression is zero as eigenvectors of an observable A corresponding to distinct eigenvalues are orthogonal. The completeness relation is merely a restatement of the resolution of the identity in terms of projectors. \square

We are now ready to do our final, complete statement of the axioms of quantum measurement.

Axiom: Quantum measurement (version 3/final)

Let $A = \sum_a a\Pi_a$ be the observable (a Hermitian operator) being measured, where the a s are the eigenvalues of A and $\{\Pi_a = \sum_{a_i=a} |a_i\rangle\langle a_i|\}_a$ are the associated projectors onto the eigenvalue- a subspaces. Let $|\psi\rangle$ be the pre-measurement state.

Dirac Postulate: If outcome a is measured, then the post measurement state is given by:

$$|\psi\rangle \mapsto \frac{1}{\sqrt{\langle\psi|\Pi_a|\psi\rangle}} \Pi_a|\psi\rangle. \quad (1.82)$$

Born rule: The probability of measuring outcome a is given by:

$$p(a) = \langle\psi|\Pi_a|\psi\rangle. \quad (1.83)$$

We have now reproduced the form of the Dirac postulate and Born rule shown in the initial table! The above formulation of the measurement axiom(s) is very general, encompassing possible degenerate eigenvalues in the measured observables. Of course, it should be consistent with our previous statement

of the axioms in the case that the eigenvalues are non-degenerate. It is highly recommended that you try this as an exercise first, but we will also give the argument here.

If an eigenvalue a is non degenerate, then $\Pi_a = |a\rangle\langle a|$. Eq. (1.82) then reads:

$$|\psi\rangle \mapsto \frac{1}{\sqrt{\langle\psi|a\rangle\langle a|\psi\rangle}} |a\rangle\langle a|\psi\rangle = \frac{\langle a|\psi\rangle}{\sqrt{|\langle a|\psi\rangle|^2}} |a\rangle = \frac{\langle a|\psi\rangle}{|\langle a|\psi\rangle|} |a\rangle = e^{i\phi} |a\rangle \sim |a\rangle. \quad (1.84)$$

Which is consistent with the previous statement of the Dirac postulate. For the Born rule, we see that Eq. (1.83) reads:

$$p(a) = \langle\psi|a\rangle\langle a|\psi\rangle = |\langle a|\psi\rangle|^2 \quad (1.85)$$

which is again consistent with our previous formulation.

The astute reader may object that we have seemingly ignored the possible complex phase factor sitting out front in Eq. (1.84). However, this is not being sloppy, but instead a fact about quantum states that *global phases are irrelevant*.

Theorem: Irrelevance of global phase

$|\psi\rangle$ and $|\varphi\rangle = e^{i\phi}|\psi\rangle$ correspond to the same physical quantum state.

Proof. The two states are only distinct if we are able to distinguish them in a measurement. However, when calculating the probability of measuring an arbitrary outcome a for any observable A with the Born rule, we find that the two have identical measurement statistics:

$$p_\varphi(a) = \langle\varphi|\Pi_a|\varphi\rangle = \langle\psi|e^{-i\phi}\Pi_a e^{i\phi}|\psi\rangle = \langle\psi|e^{-i\phi}e^{i\phi}\Pi_a|\psi\rangle = \langle\psi|\Pi_a|\psi\rangle = p_\psi(a). \quad (1.86)$$

They therefore represent the same quantum state. \square

We however note that *relative phases are significant/relevant*. For example, the S_x eigenstates $|+\rangle = \frac{|\uparrow\rangle+|\downarrow\rangle}{\sqrt{2}}$ and $|-\rangle = \frac{|\uparrow\rangle-|\downarrow\rangle}{\sqrt{2}}$ differ by a relative phase, and are hence different quantum states.

Let us return to our motivating example with the spin-1 particle. We established that:

$$S_z^2 = \hbar^2 (|+\rangle\langle +| + |-\rangle\langle -|) \quad (1.87)$$

has a degenerate eigenvalue, with both $|+\rangle$ and $|-\rangle$ being eigenstates with eigenvalue $+\hbar^2$. To deal with this degeneracy, we can use our new projector formalism of measurement. The projector corresponding to the \hbar^2 subspace is given by:

$$\Pi_{\hbar^2} = |+\rangle\langle +| + |-\rangle\langle -| \quad (1.88)$$

while the projector corresponding to the eigenvalue 0 subspace is given by:

$$\Pi_0 = |0\rangle\langle 0|. \quad (1.89)$$

So, if we wanted to find the probability of measuring $S_z^2 = \hbar^2$ given a pre-measurement state $|\psi\rangle$, the probability would be given by:

$$p(\hbar^2) = \langle\psi|\Pi_1|\psi\rangle = |\langle +|\psi\rangle|^2 + |\langle -|\psi\rangle|^2 \quad (1.90)$$

and the post measurement state would be given by:

$$|\psi\rangle \mapsto \frac{1}{\sqrt{\langle\psi|\Pi_{\hbar^2}|\psi\rangle}} \Pi_{\hbar^2}|\psi\rangle = \frac{1}{\sqrt{|\langle +|\psi\rangle|^2 + |\langle -|\psi\rangle|^2}} (\langle +|\psi\rangle|+\rangle + \langle -|\psi\rangle|-\rangle). \quad (1.91)$$

To conclude this section, we revisit the idea of individual vs. averaged measurement outcomes. We again consider the measurement of an observable $\sum_a a\Pi_a$. As before, the measurement outcomes of individual measurements are given by the eigenvalues a of A . We can again calculate the average outcome/expectation value to be:

$$\langle A \rangle_\psi := \sum_a a p(a) = \sum_a a \langle \psi | \Pi_a | \psi \rangle = \langle \psi | \left(\sum_a a \Pi_a \right) | \psi \rangle = \langle \psi | A | \psi \rangle \quad (1.92)$$

so we see that Eq. (1.48) holds as in the non-degenerate case.

1.5 Compatible and Incompatible Observables

We have seen in the previous sections how quantum measurement is an active process; through measurement the input state is projected into a subspace (corresponding to the measured eigenvalue). This has implications for whether the value of two quantum observables can be simultaneously known; we will illuminate this with a motivating example before exploring this idea more rigorously. Consider a *sequential* Stern-Gerlach experiment, as pictured in Fig. 1.7.

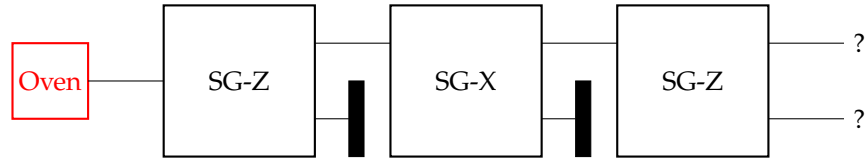


Figure 1.7: Cartoon of a sequential Stern-Gerlach experiment. First, the z -component of the spin of particles coming out of the oven area measured. The $S_z = +\frac{\hbar}{2}$ (spin up/ $|\uparrow\rangle$) particles are allowed to go through, while the $S_z = -\frac{\hbar}{2}$ (spin down/ $|\downarrow\rangle$) particles are blocked. Then, the x -component of the spin is measured, with the $S_x = +\frac{\hbar}{2}$ (spin right/ $|+\rangle$) particles allowed to go through, and the $S_x = -\frac{\hbar}{2}$ (spin left/ $|-\rangle$) particles blocked. Finally, the z -component of the spin is measured again.

Let us carry out the calculation to deduce what we would find at the output ports of the above experiment. In the first S_z measurement, there is a 50/50 probability of measuring the spin to be spin-up, and so 50% of the original particles go through. Then recalling that $|\pm\rangle = \frac{|\uparrow\rangle \pm |\downarrow\rangle}{\sqrt{2}}$, we have that $|\uparrow\rangle = \frac{|+\rangle + |-\rangle}{\sqrt{2}}$ and so we find in the second measurement that $p(+)=|\langle +|\uparrow\rangle|^2=\frac{1}{2}$ And $p(-)=|\langle -|\uparrow\rangle|^2=\frac{1}{2}$ (we may use the non-degenerate formulation of the Born rule here) so half of the particles go through - one quarter of the original particles from the oven. Now, we note something interesting; when we do the probability calculation for the third measurement, we find $p(\uparrow)=|\langle \uparrow|+\rangle|^2=\frac{1}{2}$ and $p(\downarrow)=|\langle \downarrow|+\rangle|^2=\frac{1}{2}$, so we find that one eighth of the original particles from the oven are present at each of the two output ports. Classically, this is quite strange - a classical measurement does not disturb the system, so when we measure the z -component of spin for the second time, we should expect to get 100% spin-up (as we post-selected on spin-up in the first measurement)! But this is indeed not what happens quantum mechanically - in QM, measurements physically change the state of the system, so there are physical observables that are incompatible with each other (that is, their value cannot be simultaneously precisely known). Let us formalize this notion with a definition:

Definition: Compatible Observables

Two observables A, B are *compatible* if:

$$[A, B] := AB - BA = 0. \quad (1.93)$$

The $[\cdot, \cdot]$ appearing in the above definition is known as the *commutator*, and can be interpreted as quantifying the extent to which two objects commute.

Compatible observables are characterized by the following properties:

- The probability $p(a, b)$ for obtaining the outcome a for A and b for B is independent of the order of measurement.
- Compatible measurements do not disturb one another; suppose A is measured with outcome a and then B with outcome b . If A is measured a second time, then the outcome is a again, with certainty.

These properties follow as the result of the following Lemma:

Lemma

Compatible observables have a joint eigenbasis.

Proof. Choose an (orthonormal) eigenbasis $\{|a\rangle\}$ of A . Since $[A, B] = 0$, it follows that $\langle a'|[A, B]|a\rangle = 0$ and so:

$$0 = \langle a'|[A, B]|a\rangle = (\langle a'|A)B|a\rangle - \langle a'|B(A|a\rangle) = (a' - a)\langle a'|B|a\rangle \quad (1.94)$$

So if $a \neq a'$, it follows that $\langle a'|B|a\rangle = 0$. If A is non-degenerate, we are immediately done as B is diagonal in the basis $\{|a\rangle\}$ and hence it is also an eigenbasis of B . However, if A is degenerate then $a = a'$ is possible for distinct $|a\rangle, |a'\rangle$. This means that B is block diagonal in the basis $\{|a\rangle\}$. For example in the case where A has two degenerate eigenvalues a_1 and three degenerate eigenvalues a_2 , B in this eigenbasis will be block diagonal with two blocks:

$$A = \begin{pmatrix} \boxed{\begin{matrix} a_1 & 0 \\ 0 & a_1 \end{matrix}} & 0 \\ 0 & \boxed{\begin{matrix} a_2 & 0 & 0 \\ 0 & a_2 & 0 \\ 0 & 0 & a_2 \end{matrix}} \end{pmatrix} \quad B = \begin{pmatrix} \blacksquare & 0 \\ 0 & \blacksquare \end{pmatrix} \quad (1.95)$$

Now, we borrow a known theorem from linear algebra (the Spectral theorem) that a Hermitian matrix can be diagonalized with a unitary matrix - hence, there exists some unitary transformation U (which will have a block-diagonal matrix representation consisting of blocks U_1, U_2 in the same location as the blocks of B) which diagonalizes B :

$$U = \begin{pmatrix} \boxed{U_1} & 0 \\ 0 & \boxed{U_2} \end{pmatrix} \quad (1.96)$$

Then, we have that $UAU^\dagger = A$ (consider that the blocks of U act on the eigenspaces of A independently, and then cancel to the identity by the condition $UU^\dagger = I$, thus leaving A invariant) and $UBU^\dagger = B'$ is diagonal. In other words, the basis $\{U^\dagger|a\rangle\}$ is a joint eigenbasis of A, B and we are done. \square

Since A, B share an eigenbasis, in this same basis the projectors corresponding to the various are diagonal. It therefore follows that:

$$[\Pi_{A,a}, \Pi_{B,b}] = [\Pi_{A,a}, B] = [A, \Pi_{B,b}] = 0. \quad (1.97)$$

From this fact, we can argue the two properties of compatible observables. First, we show that the probability $p(a, b)$ for obtaining the outcome a for A and b for B is independent of the order of measurement.

Suppose we measure A first, then B . Then the post measurement state after the first measurement is $\frac{\Pi_{A,a}|\psi\rangle}{\sqrt{\langle\psi|\Pi_{A,a}|\psi\rangle}}$, by the Dirac postulate. The probability to measure b in the second measurement (given that we measured a in the first measurement) is then given by the Born rule to be:

$$p(b|a) = \frac{\langle\psi|\Pi_{A,a}\Pi_{B,b}\Pi_{A,a}|\psi\rangle}{\langle\psi|\Pi_{A,a}|\psi\rangle} \quad (1.98)$$

The probability for measuring a and b is then calculated as:

$$\begin{aligned} p(a, b) &= p(a \cap b) = p(b|a)p(a) = \left(\frac{\langle\psi|\Pi_{A,a}\Pi_{B,b}\Pi_{A,a}|\psi\rangle}{\langle\psi|\Pi_{A,a}|\psi\rangle} \right) \langle\psi|\Pi_{A,a}|\psi\rangle \\ &= \langle\psi|\Pi_{A,a}\Pi_{B,b}\Pi_{A,a}|\psi\rangle \\ &= \langle\psi|\Pi_{A,a}^2\Pi_{B,b}|\psi\rangle \\ &= \langle\psi|\Pi_{A,a}\Pi_{B,b}|\psi\rangle \end{aligned} \quad (1.99)$$

where in the second-to-last equality we use that the projectors commute, and in the last equality we use that projectors are idempotent. If we measure B first and then A and go through the same calculation (simply by interchanging $A \leftrightarrow B$), we find the exact same expression, and therefore we conclude that $p(a, b)$ is independent of measurement order.

Let us also demonstrate that the compatible measurements do not disturb one another. If we measure A with outcome a , we have $|\psi\rangle \mapsto |a\rangle$ with $A|a\rangle = a|a\rangle$. If we then measure B , we have $|a\rangle \mapsto \frac{\Pi_{B,b}|a\rangle}{\sqrt{\langle a|\Pi_{B,b}|a\rangle}}$. Using the fact that $A, \Pi_{B,b}$ commute we then observe:

$$A \left(\frac{\Pi_{B,b}|a\rangle}{\sqrt{\langle a|\Pi_{B,b}|a\rangle}} \right) = \frac{\Pi_{B,b}A|a\rangle}{\sqrt{\langle a|\Pi_{B,b}|a\rangle}} = a \left(\frac{\Pi_{B,b}|a\rangle}{\sqrt{\langle a|\Pi_{B,b}|a\rangle}} \right) \quad (1.100)$$

so the post-measurement state after measuring B is still an eigenstate of A , with the same eigenvalue a !

We note an important example of compatible observables. For all observables A , it holds that $[A^n, A^m] = 0$ for all $n, m \in \mathbb{N}$. In other words, all of $\{A^k, k \in \mathbb{N}\}$ can be simultaneously measured. In the homework, you will be tasked with considering the relationship between the expectation values $|A^k|_\psi = \langle\psi|A^k|\psi\rangle$ of observables A^k , and the measurement probabilities $p(i) = |\langle b_i|\psi\rangle|^2$ where $\{|b_i\rangle\}_i$ is the eigenbasis of A .

Alongside the discussion of compatible observables runs the consideration of incompatible observables:

Definition: Incompatible Observables

Two observables A, B are *incompatible* if:

$$[A, B] \neq 0. \quad (1.101)$$

Incompatible observables have the properties that:

- The probability $p(a, b)$ for obtaining the outcome a for A and b for B is generally dependent of the order of measurement.
- Incompatible measurements disturb one another.

A good example is our motivating example of the S_z and S_x measurements at the beginning of this section. Indeed, $[S_z, S_x] = i\hbar S_y \neq 0$ and so the two observables are incompatible.

1.6 Position and Momentum

In this section, we discuss two key examples of observables - namely position and momentum. We discuss their eigenstates, which take on a continuous range. We discuss their commutation relation which is the axiomatic foundation of how they relate. We derive the momentum operator in the position basis, and show that position and momentum wavefunctions relate via a Fourier transform. Finally, we will derive the famous Heisenberg uncertainty relation.

2 Quantum Dynamics

3 Quantum Foundations

4 Symmetries and Angular Momentum

5 **Approximation Methods**

6 Identical Particles