

# Solid State Physics

**University of Cambridge Part II Mathematical Tripos**

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## Recommended Books and Resources

There are many excellent books on Solid State Physics. The two canonical books are

- Ashcroft and Mermin, *Solid State Physics*
- Kittel, *Introduction to Solid State Physics*

Both of these go substantially beyond the material covered in this course. Personally, I have a slight preference for the verbosity of Ashcroft and Mermin.

A somewhat friendlier, easier going book is

- Steve Simon, *Solid State Physics Basics*

It covers only the basics, but does so very well. (An earlier draft can be downloaded from Steve Simon's homepage; see below for a link.)

A number of lecture notes are available on the web. Links can be found on the course webpage: <http://www.damtp.cam.ac.uk/user/tong/solidstate.html>

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This material is taught as part of the “*Applications in Quantum Mechanics*” course of the Cambridge mathematical tripos.

## 0. Introduction

Solid state physics is the study of “stuff”, of how the wonderfully diverse properties of solids can emerge from the simple laws that govern electrons and atoms.

There is one, over-riding, practical reason for wanting to understand the behaviour of stuff: this is how we build things. In particular, it is how we build the delicate and powerful technologies that underlie our society. Important though they are, such practicalities will take a back seat in our story. Instead, our mantra is “knowledge for its own sake”. Indeed, the subject of solid state physics turns out to be one of extraordinary subtlety and beauty. If such knowledge ultimately proves useful, this is merely a happy corollary.

We will develop only the basics of solid state physics. We will learn how electrons glide through seemingly impenetrable solids, how their collective motion is described by a Fermi surface, and how the vibrations of the underlying atoms get tied into bundles of energy known as phonons. We will learn that electrons in magnetic fields can do strange things and start to explore some of the roles that geometry and topology play in quantum physics.

One of the ultimate surprises of solid state physics is how the subject later dovetails with ideas from particle physics. At first glance, one might have thought these two disciplines should have nothing to do with each other. Yet one of the most striking themes in modern physics is how ideas from one have influenced the other. In large part this is because both subjects rest on some of the deepest principles in physics: ideas such as symmetry, topology and universality. Although much of what we cover in these lectures will be at a basic level, we will nonetheless see some hints of these deeper connections. We will, for example, see the Dirac equation — originally introduced to unify relativity and quantum mechanics — emerging from graphene. We will learn how the vibrations of a lattice, and the resulting phonons, provide a baby introduction to quantum field theory.

# 1. Particles in a Magnetic Field

The purpose of this chapter is to understand how quantum particles react to magnetic fields. In contrast to later sections, we will not yet place these particles inside solids, for the simple reason that there is plenty of interesting behaviour to discover before we do this. Later, in Section 3.1, we will understand how these magnetic fields affect the electrons in solids.

Before we get to describe quantum effects, we first need to highlight a few of the more subtle aspects that arise when discussing classical physics in the presence of a magnetic field.

## 1.1 Gauge Fields

Recall from our lectures on [Electromagnetism](#) that the electric field  $\mathbf{E}(\mathbf{x}, t)$  and magnetic field  $\mathbf{B}(\mathbf{x}, t)$  can be written in terms a scalar potential  $\phi(\mathbf{x}, t)$  and a vector potential  $\mathbf{A}(\mathbf{x}, t)$ ,

$$\mathbf{E} = -\nabla\phi - \frac{\partial\mathbf{A}}{\partial t} \quad \text{and} \quad \mathbf{B} = \nabla \times \mathbf{A} \quad (1.1)$$

Both  $\phi$  and  $\mathbf{A}$  are referred to as *gauge fields*. When we first learn electromagnetism, they are introduced merely as handy tricks to help solve the Maxwell equations. However, as we proceed through theoretical physics, we learn that they play a more fundamental role. In particular, they are necessary if we want to discuss a Lagrangian or Hamiltonian approach to electromagnetism. We will soon see that these gauge fields are quite indispensable in quantum mechanics.

The Lagrangian for a particle of charge  $q$  and mass  $m$  moving in a background electromagnetic fields is

$$L = \frac{1}{2}m\dot{\mathbf{x}}^2 + q\dot{\mathbf{x}} \cdot \mathbf{A} - q\phi \quad (1.2)$$

The classical equation of motion arising from this Lagrangian is

$$m\ddot{\mathbf{x}} = q(\mathbf{E} + \dot{\mathbf{x}} \times \mathbf{B})$$

This is the *Lorentz force law*.

Before we proceed I should warn you of a minus sign issue. We will work with a general charge  $q$ . However, many textbooks work with the charge of the electron, written as  $q = -e$ . If this minus sign leans to confusion, you should blame Benjamin Franklin.

## An Example: Motion in a Constant Magnetic Field

We'll take a constant magnetic field, pointing in the  $z$ -direction:  $\mathbf{B} = (0, 0, B)$ . We'll take  $\mathbf{E} = 0$ . The particle is free in the  $z$ -direction, with the equation of motion  $m\ddot{z} = 0$ . The more interesting dynamics takes place in the  $(x, y)$ -plane where the equations of motion are

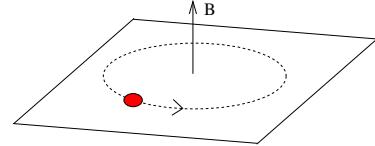
$$m\ddot{x} = qBy \quad \text{and} \quad m\ddot{y} = -qBx \quad (1.3)$$

which has general solution is

$$x(t) = X + R \sin(\omega_B(t - t_0)) \quad \text{and} \quad y(t) = Y + R \cos(\omega_B(t - t_0))$$

We see that the particle moves in a circle which, for  $B > 0$  and  $q > 0$ , is in a clockwise direction. The *cyclotron frequency* is defined by

$$\omega_B = \frac{qB}{m} \quad (1.4)$$



**Figure 1:**

The centre of the circle  $(X, Y)$ , the radius of the circle  $R$  and the phase  $t_0$  are all arbitrary. These are the four integration constants expected in the solution of two, second order differential equations.

### 1.1.1 The Hamiltonian

The *canonical momentum* in the presence of gauge fields is

$$\mathbf{p} = \frac{\partial L}{\partial \dot{\mathbf{x}}} = m\dot{\mathbf{x}} + q\mathbf{A} \quad (1.5)$$

This clearly is not the same as what we naively call momentum, namely  $m\dot{\mathbf{x}}$ .

The Hamiltonian is given by

$$H = \dot{\mathbf{x}} \cdot \mathbf{p} - L = \frac{1}{2m}(\mathbf{p} - q\mathbf{A})^2 + q\phi$$

Written in terms of the velocity of the particle, the Hamiltonian looks the same as it would in the absence of a magnetic field:  $H = \frac{1}{2}m\dot{\mathbf{x}}^2 + q\phi$ . This is the statement that a magnetic field does no work and so doesn't change the energy of the system. However, there's more to the Hamiltonian framework than just the value of  $H$ . We need to remember which variables are canonical. This information is encoded in the Poisson bracket structure of the theory (or, in fancy language, the symplectic structure on phase space). The fact that  $\mathbf{x}$  and  $\mathbf{p}$  are canonical means that

$$\{x_i, p_j\} = \delta_{ij} \quad \text{with} \quad \{x_i, x_j\} = \{p_i, p_j\} = 0$$

In the quantum theory, this structure transferred onto commutation relations between operators, which become

$$[x_i, p_j] = i\hbar\delta_{ij} \quad \text{with} \quad [x_i, x_j] = [p_i, p_j] = 0$$

### 1.1.2 Gauge Transformations

The gauge fields  $\mathbf{A}$  and  $\phi$  are not unique. We can change them as

$$\phi \rightarrow \phi - \frac{\partial \alpha}{\partial t} \quad \text{and} \quad \mathbf{A} \rightarrow \mathbf{A} + \nabla \alpha \quad (1.6)$$

for any function  $\alpha(\mathbf{x}, t)$ . Under these transformations, the electric and magnetic fields (1.1) remain unchanged. The Lagrangian (1.2) changes by a total derivative, but this is sufficient to ensure that the resulting equations of motion (1.3) are unchanged. Different choices of  $\alpha$  are said to be different choices of *gauge*. We'll see some examples below.

The existence of gauge transformations is a redundancy in our description of the system: fields which differ by the transformation (1.6) describe physically identical configurations. Nothing that we can physically measure can depend on our choice of gauge. This, it turns out, is a beautifully subtle and powerful restriction. We will start to explore some of these subtleties in Sections 1.3 and 1.4

The canonical momentum  $\mathbf{p}$  defined in (1.5) is not gauge invariant: it transforms as  $\mathbf{p} \rightarrow \mathbf{p} + q\nabla\alpha$ . This means that the numerical value of  $\mathbf{p}$  can't have any physical meaning since it depends on our choice of gauge. In contrast, the velocity of the particle  $\dot{\mathbf{x}}$  is gauge invariant, and therefore physical.

### The Schrödinger Equation

Finally, we can turn to the quantum theory. We'll look at the spectrum in the next section, but first we wish to understand how gauge transformations work. Following the usual quantisation procedure, we replace the canonical momentum with

$$\mathbf{p} \mapsto -i\hbar\nabla$$

The time-dependent Schrödinger equation for a particle in an electric and magnetic field then takes the form

$$i\hbar \frac{\partial \psi}{\partial t} = H\psi = \frac{1}{2m} \left( -i\hbar\nabla - q\mathbf{A} \right)^2 \psi + q\phi\psi \quad (1.7)$$

The shift of the kinetic term to incorporate the vector potential  $\mathbf{A}$  is sometimes referred to as *minimal coupling*.

Before we solve for the spectrum, there are two lessons to take away. The first is that it is not possible to formulate the quantum mechanics of particles moving in electric and magnetic fields in terms of  $\mathbf{E}$  and  $\mathbf{B}$  alone. We're obliged to introduce the gauge fields  $\mathbf{A}$  and  $\phi$ . This might make you wonder if, perhaps, there is more to  $\mathbf{A}$  and  $\phi$  than we first thought. We'll see the answer to this question in Section 1.3. (Spoiler: the answer is yes.)

The second lesson follows from looking at how (1.7) fares under gauge transformations. It is simple to check that the Schrödinger equation transforms covariantly (i.e. in a nice way) only if the wavefunction itself also transforms with a position-dependent phase

$$\psi(\mathbf{x}, t) \rightarrow e^{iq\alpha(\mathbf{x}, t)/\hbar} \psi(\mathbf{x}, t) \quad (1.8)$$

This is closely related to the fact that  $\mathbf{p}$  is not gauge invariant in the presence of a magnetic field. Importantly, this gauge transformation does not affect physical probabilities which are given by  $|\psi|^2$ .

The simplest way to see that the Schrödinger equation transforms nicely under the gauge transformation (1.8) is to define the *covariant derivatives*

$$\mathcal{D}_t = \frac{\partial}{\partial t} + \frac{iq}{\hbar} \phi \quad \text{and} \quad \mathcal{D}_i = \frac{\partial}{\partial x^i} - \frac{iq}{\hbar} A_i$$

In terms of these covariant derivatives, the Schrödinger equation becomes

$$i\hbar \mathcal{D}_t \psi = -\frac{\hbar^2}{2m} \mathcal{D}^2 \psi \quad (1.9)$$

But these covariant derivatives are designed to transform nicely under a gauge transformation (1.6) and (1.8). You can check that they pick up only a phase

$$\mathcal{D}_t \psi \rightarrow e^{iq\alpha/\hbar} \mathcal{D}_t \psi \quad \text{and} \quad \mathcal{D}_i \psi \rightarrow e^{iq\alpha/\hbar} \mathcal{D}_i \psi$$

This ensures that the Schrödinger equation (1.9) transforms covariantly.

## 1.2 Landau Levels

Our task now is to solve for the spectrum and wavefunctions of the Schrödinger equation. We are interested in the situation with vanishing electric field,  $\mathbf{E} = 0$ , and constant magnetic field. The quantum Hamiltonian is

$$H = \frac{1}{2m} (\mathbf{p} - q\mathbf{A})^2 \quad (1.10)$$

We take the magnetic field to lie in the  $z$ -direction, so that  $\mathbf{B} = (0, 0, B)$ . To proceed, we need to find a gauge potential  $\mathbf{A}$  which obeys  $\nabla \times \mathbf{A} = \mathbf{B}$ . There is, of course, no unique choice. Here we pick

$$\mathbf{A} = (0, xB, 0) \quad (1.11)$$

This is called *Landau gauge*. Note that the magnetic field  $\mathbf{B} = (0, 0, B)$  is invariant under both translational symmetry and rotational symmetry in the  $(x, y)$ -plane. However, the choice of  $\mathbf{A}$  is not; it breaks translational symmetry in the  $x$  direction (but not in the  $y$  direction) and rotational symmetry. This means that, while the physics will be invariant under all symmetries, the intermediate calculations will not be manifestly invariant. This kind of compromise is typical when dealing with magnetic field.

The Hamiltonian (1.10) becomes

$$H = \frac{1}{2m} (p_x^2 + (p_y - qBx)^2 + p_z^2)$$

Because we have manifest translational invariance in the  $y$  and  $z$  directions, we have  $[p_y, H] = [p_z, H] = 0$  and can look for energy eigenstates which are also eigenstates of  $p_y$  and  $p_z$ . This motivates the ansatz

$$\psi(\mathbf{x}) = e^{ik_y y + ik_z z} \chi(x) \quad (1.12)$$

Acting on this wavefunction with the momentum operators  $p_y = -i\hbar\partial_y$  and  $p_z = -i\hbar\partial_z$ , we have

$$p_y \psi = \hbar k_y \psi \quad \text{and} \quad p_z \psi = \hbar k_z \psi$$

The time-independent Schrödinger equation is  $H\psi = E\psi$ . Substituting our ansatz (1.12) simply replaces  $p_y$  and  $p_z$  with their eigenvalues, and we have

$$H\psi(\mathbf{x}) = \frac{1}{2m} \left[ p_x^2 + (\hbar k_y - qBx)^2 + \hbar^2 k_z^2 \right] \psi(\mathbf{x}) = E\psi(\mathbf{x})$$

We can write this as an eigenvalue equation for the equation  $\chi(x)$ . We have

$$\tilde{H}\chi(x) = \left( E - \frac{\hbar^2 k_z^2}{2m} \right) \chi(x)$$

where  $\tilde{H}$  is something very familiar: it's the Hamiltonian for a harmonic oscillator in the  $x$  direction, with the centre displaced from the origin,

$$\tilde{H} = \frac{1}{2m} p_x^2 + \frac{m\omega_B^2}{2} (x - k_y l_B^2)^2 \quad (1.13)$$

The frequency of the harmonic oscillator is again the cyclotron frequency  $\omega_B = qB/m$ , and we've also introduced a length scale  $l_B$ . This is a characteristic length scale which governs any quantum phenomena in a magnetic field. It is called the *magnetic length*.

$$l_B = \sqrt{\frac{\hbar}{qB}}$$

To give you some sense for this, in a magnetic field of  $B = 1$  Tesla, the magnetic length for an electron is  $l_B \approx 2.5 \times 10^{-8}$  m.

Something rather strange has happened in the Hamiltonian (1.13): the momentum in the  $y$  direction,  $\hbar k_y$ , has turned into the position of the harmonic oscillator in the  $x$  direction, which is now centred at  $x = k_y l_B^2$ .

We can immediately write down the energy eigenvalues of (1.13); they are simply those of the harmonic oscillator

$$E = \hbar\omega_B \left( n + \frac{1}{2} \right) + \frac{\hbar^2 k_z^2}{2m} \quad n = 0, 1, 2, \dots \quad (1.14)$$

The wavefunctions depend on three quantum numbers,  $n \in \mathbf{N}$  and  $k_y, k_z \in \mathbf{R}$ . They are

$$\psi_{n,k}(x, y) \sim e^{ik_y y + ik_z z} H_n(x - k_y l_B^2) e^{-(x - k_y l_B^2)^2 / 2l_B^2} \quad (1.15)$$

with  $H_n$  the usual Hermite polynomial wavefunctions of the harmonic oscillator. The  $\sim$  reflects the fact that we have made no attempt to normalise these wavefunctions.

The wavefunctions look like strips, extended in the  $y$  direction but exponentially localised around  $x = k_y l_B^2$  in the  $x$  direction. However, you shouldn't read too much into this. As we will see shortly, there is large degeneracy of wavefunctions and by taking linear combinations of these states we can cook up wavefunctions that have pretty much any shape you like.

### 1.2.1 Degeneracy

The dynamics of the particle in the  $z$ -direction is unaffected by the magnetic field  $\mathbf{B} = (0, 0, B)$ . To focus on the novel physics, let's restrict to particles with  $k_z = 0$ . The energy spectrum then coincides with that of a harmonic oscillator,

$$E_n = \hbar\omega_B \left( n + \frac{1}{2} \right) \quad (1.16)$$

In the present context, these are called *Landau levels*. We see that, in the presence of a magnetic field, the energy levels of a particle become equally spaced, with the gap between each level proportional to the magnetic field  $B$ . Note that the energy spectrum looks very different from a free particle moving in the  $(x, y)$ -plane.

The states in a given Landau level are not unique. Instead, there is a huge degeneracy, with many states having the same energy. We can see this in the form of the wavefunctions (1.15) which, when  $k_z = 0$ , depend on two quantum numbers,  $n$  and  $k_y$ . Yet the energy (1.16) is independent of  $k_y$ .

Let's determine how large this degeneracy of states is. To do so, we need to restrict ourselves to a finite region of the  $(x, y)$ -plane. We pick a rectangle with sides of lengths  $L_x$  and  $L_y$ . We want to know how many states fit inside this rectangle.

Having a finite size  $L_y$  is like putting the system in a box in the  $y$ -direction. The wavefunctions must obey

$$\psi(x, y + L_y, z) = \psi(x, y, z) \quad \Rightarrow \quad e^{ik_y L_y} = 1$$

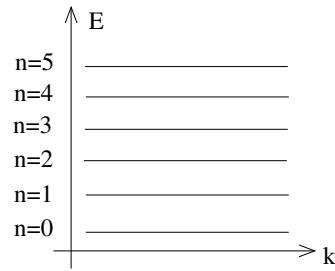
This means that the momentum  $k_y$  is quantised in units of  $2\pi/L_y$ .

Having a finite size  $L_x$  is somewhat more subtle. The reason is that, as we mentioned above, the gauge choice (1.11) does not have manifest translational invariance in the  $x$ -direction. This means that our argument will be a little heuristic. Because the wavefunctions (1.15) are exponentially localised around  $x = k_y l_B^2$ , for a finite sample restricted to  $0 \leq x \leq L_x$  we would expect the allowed  $k_y$  values to range between  $0 \leq k_y \leq L_x/l_B^2$ . The end result is that the number of states in each Landau level is given by

$$\mathcal{N} = \frac{L_y}{2\pi} \int_0^{L_x/l_B^2} dk = \frac{L_x L_y}{2\pi l_B^2} = \frac{qBA}{2\pi\hbar} \quad (1.17)$$

where  $A = L_x L_y$  is the area of the sample. Strictly speaking, we should take the integer part of the answer above.

The degeneracy (1.17) is very very large. Throwing in some numbers, there are around  $10^{10}$  degenerate states per Landau level for electrons in a region of area  $A = 1 \text{ cm}^2$  in a magnetic field  $B \sim 0.1 \text{ T}$ . This large degeneracy ultimately, this leads to an array of dramatic and surprising physics.



**Figure 2:** Landau Levels

### 1.2.2 Symmetric Gauge

It is worthwhile to repeat the calculations above using a different gauge choice. This will give us a slightly different perspective on the physics. A natural choice is *symmetric gauge*

$$\mathbf{A} = -\frac{1}{2}\mathbf{x} \times \mathbf{B} = \frac{B}{2}(-y, x, 0) \quad (1.18)$$

This choice of gauge breaks translational symmetry in both the  $x$  and the  $y$  directions. However, it does preserve rotational symmetry about the origin. This means that angular momentum is now a good quantum number to label states.

In this gauge, the Hamiltonian is given by

$$\begin{aligned} H &= \frac{1}{2m} \left[ \left( p_x + \frac{qBy}{2} \right)^2 + \left( p_y - \frac{qBx}{2} \right)^2 + p_z^2 \right] \\ &= -\frac{\hbar^2}{2m} \nabla^2 - \frac{qB}{2m} L_z + \frac{q^2 B^2}{8m} (x^2 + y^2) \end{aligned} \quad (1.19)$$

where we've introduced the angular momentum operator

$$L_z = x p_y - y p_x$$

We'll again restrict to motion in the  $(x, y)$ -plane, so we focus on states with  $k_z = 0$ . It turns out that complex variables are particularly well suited to describing states in symmetric gauge, in particular in the lowest Landau level with  $n = 0$ . We define

$$w = x + iy \quad \text{and} \quad \bar{w} = x - iy$$

Correspondingly, the complex derivatives are

$$\partial = \frac{1}{2} \left( \frac{\partial}{\partial x} - i \frac{\partial}{\partial y} \right) \quad \text{and} \quad \bar{\partial} = \frac{1}{2} \left( \frac{\partial}{\partial x} + i \frac{\partial}{\partial y} \right)$$

which obey  $\partial w = \bar{\partial} \bar{w} = 1$  and  $\partial \bar{w} = \bar{\partial} w = 0$ . The Hamiltonian, restricted to states with  $k_z = 0$ , is then given by

$$H = -\frac{2\hbar^2}{m} \partial \bar{\partial} - \frac{\omega_B}{2} L_z + \frac{m\omega_B^2}{8} w \bar{w}$$

where now

$$L_z = \hbar(w\partial - \bar{w}\bar{\partial})$$

It is simple to check that the states in the lowest Landau level take the form

$$\psi_0(w, \bar{w}) = f(w) e^{-|w|^2/4l_B^2}$$

for *any* holomorphic function  $f(w)$ . These all obey

$$H\psi_0(w, \bar{w}) = \frac{\hbar\omega_B}{2}\psi_0(w, \bar{w})$$

which is the statement that they lie in the lowest Landau level with  $n = 0$ . We can further distinguish these states by requiring that they are also eigenvalues of  $L_z$ . These are satisfied by the monomials,

$$\psi_0 = w^M e^{-|w|^2/4l_B^2} \Rightarrow L_z\psi_0 = \hbar M\psi_0 \quad (1.20)$$

for some positive integer  $M$ .

### Degeneracy Revisited

In symmetric gauge, the profiles of the wavefunctions (1.20) form concentric rings around the origin. The higher the angular momentum  $M$ , the further out the ring. This, of course, is very different from the strip-like wavefunctions that we saw in Landau gauge (1.15). You shouldn't read too much into this other than the fact that the profile of the wavefunctions is not telling us anything physical as it is not gauge invariant.

However, it's worth revisiting the degeneracy of states in symmetric gauge. The probability for a particle with angular momentum  $M$  is peaked on a ring of radius  $r = \sqrt{2M}l_B$ . This means that in a disc shaped region of area  $A = \pi R^2$ , the number of states is roughly (the integer part of)

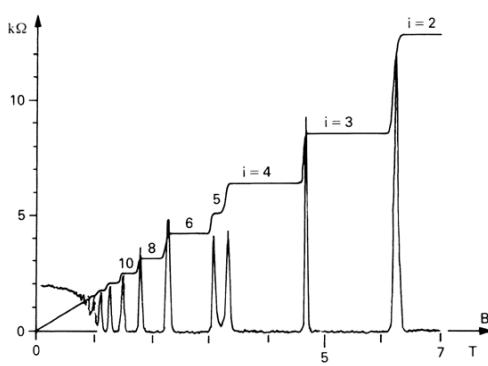
$$\mathcal{N} = R^2/2l_B^2 = A/2\pi l_B^2 = \frac{qBA}{2\pi\hbar}$$

which agrees with our earlier result (1.17).

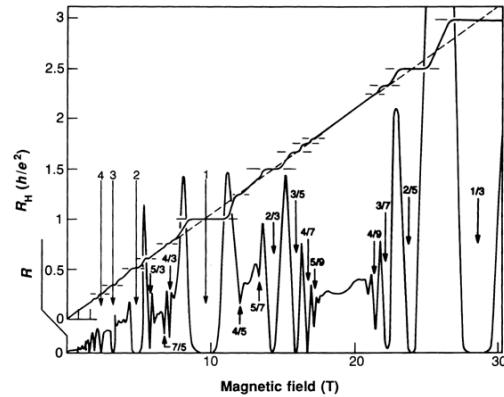
### 1.2.3 An Invitation to the Quantum Hall Effect

Take a system with some fixed number of electrons, which are restricted to move in the  $(x, y)$ -plane. The charge of the electron is  $q = -e$ . In the presence of a magnetic field, these will first fill up the  $\mathcal{N} = eBA/2\pi\hbar$  states in the  $n = 0$  lowest Landau level. If any are left over they will then start to fill up the  $n = 1$  Landau level, and so on.

Now suppose that we increase the magnetic field  $B$ . The number of states  $\mathcal{N}$  housed in each Landau level will increase, leading to a depletion of the higher Landau levels. At certain, very special values of  $B$ , we will find some number of Landau levels that are exactly filled. However, generically there will be a highest Landau level which is only partially filled.



**Figure 3:** The integer quantum Hall effect.



**Figure 4:** The fractional quantum Hall effect.

This successive depletion of Landau levels gives rise to a number of striking signatures in different physical quantities. Often these quantities oscillate, or jump discontinuously as the number of occupied Landau levels varies. One particular example is the de Haas van Alphen oscillations seen in the magnetic susceptibility which we describe in Section 3.3.4. Another example is the behaviour of the resistivity  $\rho$ . This relates the current density  $\mathbf{J} = (J_x, J_y)$  to the applied electric field  $\mathbf{E} = (E_x, E_y)$ ,

$$\mathbf{E} = \rho \mathbf{J}$$

In the presence of an applied magnetic field  $\mathbf{B} = (0, 0, B)$ , the electrons move in circles. This results in components of the current which are both parallel and perpendicular to the electric field. This is modelled straightforwardly by taking  $\rho$  to be a matrix

$$\rho = \begin{pmatrix} \rho_{xx} & \rho_{xy} \\ -\rho_{xy} & \rho_{xx} \end{pmatrix}$$

where the form of the matrix follows from rotational invariance. Here  $\rho_{xx}$  is called the *longitudinal resistivity* while  $\rho_{xy}$  is called the *Hall resistivity*.

In very clean samples, in strong magnetic fields, both components of the resistivity exhibit very surprising behaviour. This is shown in the left-hand figure above. The Hall resistivity  $\rho_{xy}$  increases with  $B$  by forming a series of plateaux, on which it takes values

$$\rho_{xy} = \frac{2\pi\hbar}{e^2} \frac{1}{\nu} \quad \nu \in \mathbf{N}$$

The value of  $\nu$  (which is labelled  $i = 2, 3, \dots$  in the data shown above) is measured to be an integer to extraordinary accuracy — around one part in  $10^9$ . Meanwhile,

the longitudinal resistivity vanishes when  $\rho_{xy}$  lies on a plateaux, but spikes whenever there is a transition between different plateaux. This phenomenon, called the *integer Quantum Hall Effect*, was discovered by Klaus von Klitzing in 1980. For this, he was awarded the Nobel prize in 1985.

It turns out that the integer quantum Hall effect is a direct consequence of the existence of discrete Landau levels. The plateaux occur when precisely  $\nu \in \mathbf{Z}^+$  Landau levels are filled. Of course, we're very used to seeing integers arising in quantum mechanics — this, after all, is what the “quantum” in quantum mechanics means. However, the quantisation of the resistivity  $\rho_{xy}$  is something of a surprise because this is a macroscopic quantity, involving the collective behaviour of many trillions of electrons, swarming through a hot and dirty system. A full understanding of the integer quantum Hall effect requires an appreciation of how the mathematics of topology fits in with quantum mechanics. David Thouless (and, to some extent, Duncan Haldane) were awarded the 2016 Nobel prize for understanding the underlying role of topology in this system.

Subsequently it was realised that similar behaviour also happens when Landau levels are partially filled. However, it doesn't occur for any filling, but only very special values. This is referred to as the *fractional quantum Hall effect*. The data is shown in the right-hand figure. You can see clear plateaux when the lowest Landau level has  $\nu = \frac{1}{3}$  of its states filled. There is another plateaux when  $\nu = \frac{2}{5}$  of the states are filled, followed by a bewildering pattern of further plateaux, all of which occur when  $\nu$  is some rational number. This was discovered by Tsui and Störmer in 1982. It called the *Fractional Quantum Hall Effect*. The 1998 Nobel prize was awarded to Tsui and Störmer, together with Laughlin who pioneered the first theoretical ideas to explain this behaviour.

The fractional quantum Hall effect cannot be explained by treating the electrons as free. Instead, it requires us to take interactions into account. We have seen that each Landau level has a macroscopically large degeneracy. This degeneracy is lifted by interactions, resulting in a new form of quantum liquid which exhibits some magical properties. For example, in this state of matter the electron — which, of course, is an indivisible particle — can split into constituent parts! The  $\nu = \frac{1}{3}$  state has excitations which carry 1/3 of the charge of an electron. In other quantum Hall states, the excitations have charge 1/5 or 1/4 of the electron. These particles also have a number of other, even stranger properties to do with their quantum statistics and there is hope that these may underly the construction of a quantum computer.

We will not delve into any further details of the quantum Hall effect. Suffice to say that it is one of the richest and most beautiful subjects in theoretical physics. You can find a fuller exploration of these ideas in the lecture notes devoted to the [Quantum Hall Effect](#).

### 1.3 The Aharonov-Bohm Effect

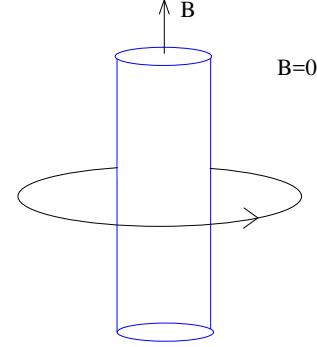
In our course on [Electromagnetism](#), we learned that the gauge potential  $A_\mu$  is unphysical: the physical quantities that affect the motion of a particle are the electric and magnetic fields. Yet we've seen above that we cannot formulate quantum mechanics without introducing the gauge fields  $\mathbf{A}$  and  $\phi$ . This might lead us to wonder whether there is more to life than  $\mathbf{E}$  and  $\mathbf{B}$  alone. In this section we will see that things are, indeed, somewhat more subtle.

#### 1.3.1 Particles Moving around a Flux Tube

Consider the set-up shown in the figure. We have a solenoid of area  $A$ , carrying magnetic field  $\mathbf{B} = (0, 0, B)$  and therefore magnetic flux  $\Phi = BA$ . Outside the solenoid the magnetic field is zero. However, the vector potential is not. This follows from Stokes' theorem which tells us that the line integral outside the solenoid is given by

$$\oint \mathbf{A} \cdot d\mathbf{x} = \int \mathbf{B} \cdot d\mathbf{S} = \Phi$$

This is simply solved in cylindrical polar coordinates by



**Figure 5:**

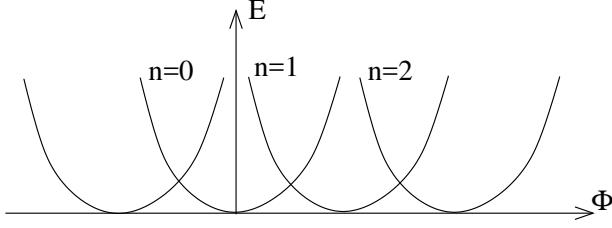
$$A_\phi = \frac{\Phi}{2\pi r}$$

Now consider a charged quantum particle restricted to lie in a ring of radius  $r$  outside the solenoid. The only dynamical degree of freedom is the angular coordinate  $\phi \in [0, 2\pi)$ . The Hamiltonian is

$$H = \frac{1}{2m} (p_\phi - qA_\phi)^2 = \frac{1}{2mr^2} \left( -i\hbar \frac{\partial}{\partial \phi} - \frac{q\Phi}{2\pi} \right)^2$$

We'd like to see how the presence of this solenoid affects the particle. The energy eigenstates are simply

$$\psi = \frac{1}{\sqrt{2\pi r}} e^{in\phi} \quad n \in \mathbf{Z} \quad (1.21)$$



**Figure 6:** The energy spectrum for a particle moving around a solenoid.

where the requirement that  $\psi$  is single valued around the circle means that we must take  $n \in \mathbf{Z}$ . Plugging this into the time independent Schrödinger equation  $H\psi = E\psi$ , we find the spectrum

$$E = \frac{1}{2mr^2} \left( \hbar n - \frac{q\Phi}{2\pi} \right)^2 = \frac{\hbar^2}{2mr^2} \left( n - \frac{\Phi}{\Phi_0} \right)^2 \quad n \in \mathbf{Z}$$

where we've defined the *quantum of flux*  $\Phi_0 = 2\pi\hbar/q$ . (Usually this quantum of flux is defined using the electron charge  $q = -e$ , with the minus signs massaged so that  $\Phi_0 \equiv 2\pi\hbar/e > 0$ .)

Note that if  $\Phi$  is an integer multiple of  $\Phi_0$ , then the spectrum is unaffected by the solenoid. But if the flux in the solenoid is not an integral multiple of  $\Phi_0$  — and there is no reason that it should be — then the spectrum gets shifted. We see that the energy of the particle knows about the flux  $\Phi$  even though the particle never goes near the region with magnetic field. The resulting energy spectrum is shown in Figure 6.

There is a slightly different way of looking at this result. Away from the solenoid, the gauge field is a total divergence

$$\mathbf{A} = \nabla\alpha \quad \text{with} \quad \alpha = \frac{\Phi\phi}{2\pi}$$

This means that we can try to remove it by redefining the wavefunction to be

$$\psi \rightarrow \tilde{\psi} = \exp\left(\frac{-iq\alpha}{\hbar}\right) \psi = \exp\left(\frac{-iq\Phi}{2\pi\hbar}\phi\right) \psi$$

However, there is an issue: the wavefunction should be single-valued. This, after all, is how we got the quantisation condition  $n \in \mathbf{Z}$  in (1.21). This means that the gauge transformation above is allowed only if  $\Phi$  is an integer multiple of  $\Phi_0 = 2\pi\hbar/q$ . Only in this case is the particle unaffected by the solenoid. The obstacle arises from the fact that the wavefunction of the particle winds around the solenoid. We see here the first glimpses of how topology starts to feed into quantum mechanics.

There are a number of further lessons lurking in this simple quantum mechanical set-up. You can read about them in the [lectures on the Quantum Hall Effect](#) (see Section 1.5.3) and the [lectures on Gauge Theory](#) (see Section 3.6.1).

### 1.3.2 Aharonov-Bohm Scattering

The fact that a quantum particle can be affected by  $\mathbf{A}$  even when restricted to regions where  $\mathbf{B} = 0$  was first pointed out by Aharonov and Bohm in a context which is closely related to the story above. They revisited the famous double-slit experiment, but now with a twist: a solenoid carrying flux  $\Phi$  is hidden behind the wall. This set-up is shown in the figure below. Once again, the particle is forbidden from going near the solenoid. Nonetheless, the presence of the magnetic flux affects the resulting interference pattern, shown as the dotted line in the figure.

Consider a particle that obeys the free Schrödinger equation,

$$\frac{1}{2m} \left( -i\hbar\nabla - q\mathbf{A} \right)^2 \psi = E\psi$$

We can formally remove the gauge field by writing

$$\psi(\mathbf{x}) = \exp \left( \frac{iq}{\hbar} \int^{\mathbf{x}} \mathbf{A}(\mathbf{x}') \cdot d\mathbf{x}' \right) \phi(\mathbf{x})$$

where the integral is over any path. Crucially, however, in the double-slit experiment there are two paths,  $P_1$  and  $P_2$ . The phase picked up by the particle due to the gauge field differs depending on which path is taken. The phase difference is given by

$$\Delta\theta = \frac{q}{\hbar} \int_{P_1} \mathbf{A} \cdot d\mathbf{x} - \frac{q}{\hbar} \int_{P_2} \mathbf{A} \cdot d\mathbf{x} = \frac{q}{\hbar} \oint \mathbf{A} \cdot d\mathbf{x} = \frac{q}{\hbar} \int \mathbf{B} \cdot d\mathbf{S}$$

Note that neither the phase arising from path  $P_1$ , nor the phase arising from path  $P_2$ , is gauge invariant. However, the difference between the two phases is gauge invariant. As we see above, it is given by the flux through the solenoid. This is the Aharonov-Bohm phase,  $e^{iq\Phi/\hbar}$ , an extra contribution that arises when charged particles move around magnetic fields.

The Aharonov-Bohm phase manifests in the interference pattern seen on the screen. As  $\Phi$  is changed, the interference pattern shifts, an effect which has been experimentally observed. Only when  $\Phi$  is an integer multiple of  $\Phi_0$  is the particle unaware of the presence of the solenoid.

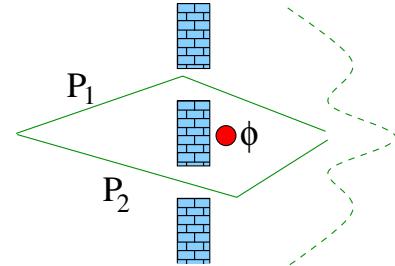


Figure 7:

## 1.4 Magnetic Monopoles

A *magnetic monopole* is a hypothetical object which emits a radial magnetic field of the form

$$\mathbf{B} = \frac{g\hat{\mathbf{r}}}{4\pi r^2} \quad \Rightarrow \quad \int d\mathbf{S} \cdot \mathbf{B} = g \quad (1.22)$$

Here  $g$  is called the *magnetic charge*.

We learned in our first course on [Electromagnetism](#) that magnetic monopoles don't exist. First, and most importantly, they have never been observed. Second there's a law of physics which insists that they can't exist. This is the Maxwell equation

$$\nabla \cdot \mathbf{B} = 0$$

Third, this particular Maxwell equation would appear to be non-negotiable. This is because it follows from the definition of the magnetic field in terms of the gauge field

$$\mathbf{B} = \nabla \times \mathbf{A} \quad \Rightarrow \quad \nabla \cdot \mathbf{B} = 0$$

Moreover, as we've seen above, the gauge field  $\mathbf{A}$  is necessary to describe the quantum physics of particles moving in magnetic fields. Indeed, the Aharonov-Bohm effect tells us that there is non-local information stored in  $\mathbf{A}$  that can only be detected by particles undergoing closed loops. All of this points to the fact that we would be wasting our time discussing magnetic monopoles any further.

Happily, there is a glorious loophole in all of these arguments, first discovered by Dirac, and magnetic monopoles play a crucial role in our understanding of the more subtle effects in gauge theories. The essence of this loophole is that there is an ambiguity in how we define the gauge potentials. In this section, we will see how this arises.

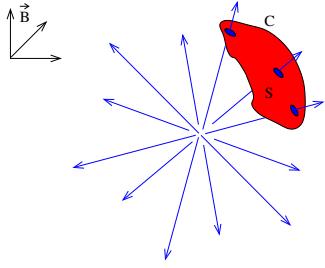
### 1.4.1 Dirac Quantisation

It turns out that not any magnetic charge  $g$  is compatible with quantum mechanics. Here we present several different arguments for the allowed values of  $g$ .

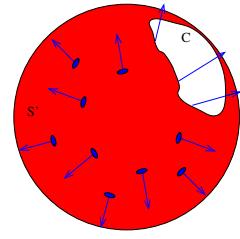
We start with the simplest and most physical of these arguments. Suppose that a particle with charge  $q$  moves along some closed path  $C$  in the background of some gauge potential  $\mathbf{A}(\mathbf{x})$ . Then, upon returning to its initial starting position, the wavefunction of the particle picks up a phase

$$\psi \rightarrow e^{iq\alpha/\hbar}\psi \quad \text{with} \quad \alpha = \oint_C \mathbf{A} \cdot d\mathbf{x} \quad (1.23)$$

This is the Aharonov-Bohm phase described above.



**Figure 8:** Integrating over  $S$ ...



**Figure 9:** ...or over  $S'$ .

The phase of the wavefunction is not an observable quantity in quantum mechanics. However, as we described above, the phase in (1.23) is really a *phase difference*. We could, for example, place a particle in a superposition of two states, one of which stays still while the other travels around the loop  $C$ . The subsequent interference will depend on the phase  $e^{iq\alpha/\hbar}$ , just like in the Aharonov-Bohm effect.

Let's now see what this has to do with magnetic monopoles. We place our particle, with electric charge  $q$ , in the background of a magnetic monopole with magnetic charge  $g$ . We keep the magnetic monopole fixed, and let the electric particle undergo some journey along a path  $C$ . We will ask only that the path  $C$  avoids the origin where the magnetic monopole is sitting. This is shown in the left-hand panel of the figure. Upon returning, the particle picks up a phase  $e^{iq\alpha/\hbar}$  with

$$\alpha = \oint_C \mathbf{A} \cdot d\mathbf{x} = \int_S \mathbf{B} \cdot d\mathbf{S}$$

where, as shown in the figure,  $S$  is the area enclosed by  $C$ . Using the fact that  $\int_{S^2} \mathbf{B} \cdot d\mathbf{S} = g$ , if the surface  $S$  makes a solid angle  $\Omega$ , this phase can be written as

$$\alpha = \frac{\Omega g}{4\pi}$$

However, there's an ambiguity in this computation. Instead of integrating over  $S$ , it is equally valid to calculate the phase by integrating over  $S'$ , shown in the right-hand panel of the figure. The solid angle formed by  $S'$  is  $\Omega' = 4\pi - \Omega$ . The phase is then given by

$$\alpha' = -\frac{(4\pi - \Omega)g}{4\pi}$$

where the overall minus sign comes because the surface  $S'$  has the opposite orientation to  $S$ . As we mentioned above, the phase shift that we get in these calculations is

observable: we can't tolerate different answers from different calculations. This means that we must have  $e^{iq\alpha/\hbar} = e^{iq\alpha'/\hbar}$ . This gives the condition

$$qg = 2\pi\hbar n \quad \text{with } n \in \mathbf{Z} \quad (1.24)$$

This is the famous Dirac quantisation condition. The smallest such magnetic charge has  $n = 1$ . It coincides with the quantum of flux,  $g = \Phi_0 = 2\pi\hbar/q$ .

Above we worked with a single particle of charge  $q$ . Obviously, the same argument must hold for any other particle of charge  $q'$ . There are two possibilities. The first is that all particles carry charge that is an integer multiple of some smallest unit. In this case, it's sufficient to impose the Dirac quantisation condition (1.24) where  $q$  is the smallest unit of charge. For example, in our world we should take  $q = \pm e$  to be the electron or proton charge (or, if we look more closely in the Standard Model, we might choose to take  $q = -e/3$ , the charge of the down quark).

The second possibility is that the particles carry electric charges which are irrational multiples of each other. For example, there may be a particle with charge  $q$  and another particle with charge  $\sqrt{2}q$ . In this case, no magnetic monopoles are allowed.

It's sometimes said that the existence of a magnetic monopole would imply the quantisation of electric charges. This, however, has it backwards. (It also misses the point that we have a wonderful explanation of the quantisation of charges from the story of anomaly cancellation in the Standard Model.) There are two possible groups that could underly gauge transformations in electromagnetism. The first is  $U(1)$ ; this has integer valued charges and admits magnetic monopoles. The second possibility is  $\mathbf{R}$ ; this has irrational electric charges and forbids monopoles. All the evidence in our world points to the fact that electromagnetism is governed by  $U(1)$  and that magnetic monopoles should exist.

Above we looked at an electrically charged particle moving in the background of a magnetically charged particle. It is simple to generalise the discussion to particles that carry both electric and magnetic charges. These are called *dyons*. For two dyons, with charges  $(q_1, g_1)$  and  $(q_2, g_2)$ , the generalisation of the Dirac quantisation condition requires

$$q_1g_2 - q_2g_1 \in 2\pi\hbar\mathbf{Z}$$

This is sometimes called the *Dirac-Zwaniger* condition.

### 1.4.2 A Patchwork of Gauge Fields

The discussion above shows how quantum mechanics constrains the allowed values of magnetic charge. It did not, however, address the main obstacle to constructing a magnetic monopole out of gauge fields  $\mathbf{A}$  when the condition  $\mathbf{B} = \nabla \times \mathbf{A}$  would seem to explicitly forbid such objects.

Let's see how to do this. Our goal is to write down a configuration of gauge fields which give rise to the magnetic field (1.22) of a monopole which we will place at the origin. However, we will need to be careful about what we want such a gauge field to look like.

The first point is that we won't insist that the gauge field is well defined at the origin. After all, the gauge fields arising from an electron are not well defined at the position of an electron and it would be churlish to require more from a monopole. This fact gives us our first bit of leeway, because now we need to write down gauge fields on  $\mathbf{R}^3/\{0\}$ , as opposed to  $\mathbf{R}^3$  and the space with a point cut out enjoys some non-trivial topology that we will make use of.

Consider the following gauge connection, written in spherical polar coordinates

$$A_\phi^N = \frac{g}{4\pi r} \frac{1 - \cos \theta}{\sin \theta} \quad (1.25)$$

The resulting magnetic field is

$$\mathbf{B} = \nabla \times \mathbf{A} = \frac{1}{r \sin \theta} \frac{\partial}{\partial \theta} (A_\phi^N \sin \theta) \hat{\mathbf{r}} - \frac{1}{r} \frac{\partial}{\partial r} (r A_\phi^N) \hat{\boldsymbol{\theta}}$$

Substituting in (1.25) gives

$$\mathbf{B} = \frac{g \hat{\mathbf{r}}}{4\pi r^2} \quad (1.26)$$

In other words, this gauge field results in the magnetic monopole. But how is this possible? Didn't we learn in kindergarten that if we can write  $\mathbf{B} = \nabla \times \mathbf{A}$  then  $\int d\mathbf{S} \cdot \mathbf{B} = 0$ ? How does the gauge potential (1.25) manage to avoid this conclusion?

The answer is that  $\mathbf{A}^N$  in (1.25) is actually a singular gauge connection. It's not just singular at the origin, where we've agreed this is allowed, but it is singular along an entire half-line that extends from the origin to infinity. This is due to the  $1/\sin \theta$  term which diverges at  $\theta = 0$  and  $\theta = \pi$ . However, the numerator  $1 - \cos \theta$  has a zero when  $\theta = 0$  and the gauge connection is fine there. But the singularity along the half-line  $\theta = \pi$  remains. The upshot is that this gauge connection is not acceptable along the line of the south pole, but is fine elsewhere. This is what the superscript  $N$  is there to remind us: we can work with this gauge connection as long as we keep north.

Now consider a different gauge connection

$$A_\phi^S = -\frac{g}{4\pi r} \frac{1 + \cos \theta}{\sin \theta} \quad (1.27)$$

This again gives rise to the magnetic field (1.26). This time it is well behaved at  $\theta = \pi$ , but singular at the north pole  $\theta = 0$ . The superscript  $S$  is there to remind us that this connection is fine as long as we keep south.

At this point, we make use of the ambiguity in the gauge connection. We are going to take  $\mathbf{A}^N$  in the northern hemisphere and  $\mathbf{A}^S$  in the southern hemisphere. This is allowed because the two gauge potentials are the same up to a gauge transformation,  $\mathbf{A} \rightarrow \mathbf{A} + \nabla \alpha$ . Recalling the expression for  $\nabla \alpha$  in spherical polars, we find that for  $\theta \neq 0, \pi$ , we can indeed relate  $A_\phi^N$  and  $A_\phi^S$  by a gauge transformation,

$$A_\phi^N = A_\phi^S + \frac{1}{r \sin \theta} \partial_\phi \alpha \quad \text{where } \alpha = \frac{g\phi}{2\pi} \quad (1.28)$$

However, there's still a question remaining: is this gauge transformation allowed? The problem is that the function  $\alpha$  is not single valued:  $\alpha(\phi = 2\pi) = \alpha(\phi = 0) + g$ . And this should concern us because, as we've seen in (1.8), the gauge transformation also acts on the wavefunction of a quantum particle

$$\psi \rightarrow e^{iq\alpha/\hbar} \psi$$

There's no reason that we should require the gauge transformation  $\alpha$  to be single-valued, but we do want the wavefunction  $\psi$  to be single-valued. This holds for the gauge transformation (1.28) provided that we have

$$qg = 2\pi\hbar n \quad \text{with } n \in \mathbf{Z}$$

This, of course, is the Dirac quantisation condition (1.24).

Mathematically, we have constructed of a topologically non-trivial  $U(1)$  bundle over the  $\mathbf{S}^2$  surrounding the origin. In this context, the integer  $n$  is called the first Chern number.

### 1.4.3 Monopoles and Angular Momentum

Here we provide yet another derivation of the Dirac quantisation condition, this time due to Saha. The key idea is that the quantisation of magnetic charge actually follows from the more familiar quantisation of angular momentum. The twist is that, in the presence of a magnetic monopole, angular momentum isn't quite what you thought.

To set the scene, let's go back to the Lorentz force law

$$\frac{d\mathbf{p}}{dt} = q\dot{\mathbf{x}} \times \mathbf{B}$$

with  $\mathbf{p} = m\dot{\mathbf{x}}$ . Recall from our discussion in Section 1.1.1 that  $\mathbf{p}$  defined here is not the canonical momentum, a fact which is hiding in the background in the following derivation. Now let's consider this equation in the presence of a magnetic monopole, with

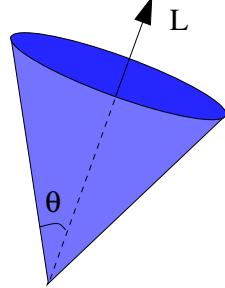
$$\mathbf{B} = \frac{q}{4\pi r^3} \frac{\mathbf{r}}{r^3}$$

The monopole has rotational symmetry so we would expect that the angular momentum,  $\mathbf{x} \times \mathbf{p}$ , is conserved. Let's check:

$$\begin{aligned} \frac{d(\mathbf{x} \times \mathbf{p})}{dt} &= \dot{\mathbf{x}} \times \mathbf{p} + \mathbf{x} \times \dot{\mathbf{p}} = \mathbf{x} \times \dot{\mathbf{p}} = q\mathbf{x} \times (\dot{\mathbf{x}} \times \mathbf{B}) \\ &= \frac{qg}{4\pi r^3} \mathbf{x} \times (\dot{\mathbf{x}} \times \mathbf{x}) = \frac{qg}{4\pi} \left( \frac{\dot{\mathbf{x}}}{r} - \frac{\dot{r}\mathbf{x}}{r^2} \right) \\ &= \frac{d}{dt} \left( \frac{qg}{4\pi} \hat{\mathbf{r}} \right) \end{aligned}$$

We see that in the presence of a magnetic monopole, the naive angular momentum  $\mathbf{x} \times \mathbf{p}$  is not conserved! However, as we also noticed in the lectures on [Classical Dynamics](#) (see Section 4.3.2), we can easily write down a modified angular momentum that is conserved, namely

$$\mathbf{L} = \mathbf{x} \times \mathbf{p} - \frac{qg}{4\pi} \hat{\mathbf{r}}$$



The extra term can be thought of as the angular momentum stored in  $\mathbf{E} \times \mathbf{B}$ . The surprise is that the system has angular momentum even when the particle doesn't move.

Before we move on, there's a nice and quick corollary that we can draw from this. The angular momentum vector  $\mathbf{L}$  does not change with time. But the angle that the particle makes with this vector is

$$\mathbf{L} \cdot \hat{\mathbf{r}} = -\frac{qg}{4\pi} = \text{constant}$$

This means that the particle moves on a cone, with axis  $\mathbf{L}$  and angle  $\cos \theta = -qg/4\pi L$ .

So far, our discussion has been classical. Now we invoke some simple quantum mechanics: the angular momentum should be quantised. In particular, the angular momentum in the  $z$ -direction should be  $L_z \in \frac{1}{2}\hbar\mathbf{Z}$ . Using the result above, we have

$$\frac{qg}{4\pi} = \frac{1}{2}\hbar n \quad \Rightarrow \quad qg = 2\pi\hbar n \quad \text{with } n \in \mathbf{Z}$$

Once again, we find the Dirac quantisation condition.

## 1.5 Spin in a Magnetic Field

As we've seen in previous courses, particles often carry an intrinsic angular momentum called *spin*  $\mathbf{S}$ . This spin is quantised in half-integer units. For examples, electrons have spin  $\frac{1}{2}$  and their spin operator is written in terms of the Pauli matrices  $\boldsymbol{\sigma}$ ,

$$\mathbf{S} = \frac{\hbar}{2}\boldsymbol{\sigma}$$

Importantly, the spin of any particle couples to a background magnetic field  $\mathbf{B}$ . The key idea here is that the intrinsic spin acts like a magnetic moment  $\mathbf{m}$  which couples to the magnetic field through the Hamiltonian

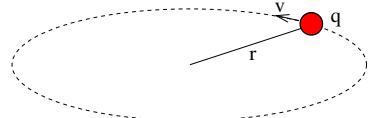
$$H = -\mathbf{m} \cdot \mathbf{B}$$

The question we would like to answer is: what magnetic moment  $\mathbf{m}$  should we associate with spin?

A full answer to this question would require an extended detour into the Dirac equation. Here we provide only some basic motivation. First consider a particle of charge  $q$  moving with velocity  $\mathbf{v}$  around a circle of radius  $\mathbf{r}$  as shown in the figure. From our lectures on [Electromagnetism](#), we know that the associated magnetic moment is given by

$$\mathbf{m} = \frac{q}{2}\mathbf{r} \times \mathbf{v} = \frac{q}{2m}\mathbf{L}$$

where  $\mathbf{L} = m\mathbf{r} \times \mathbf{v}$  is the orbital angular momentum of the particle. Indeed, we already saw the resulting coupling  $H = -(q/2m)\mathbf{L} \cdot \mathbf{B}$  in our derivation of the Hamiltonian in symmetric gauge (1.19).



**Figure 11:**

Since the spin of a particle is another contribution to the angular momentum, we might anticipate that the associated magnetic moment takes the form

$$\mathbf{m} = g \frac{q}{2m} \mathbf{S}$$

where  $g$  is some dimensionless number. (Note:  $g$  is unrelated to the magnetic charge that we discussed in the previous section!) This, it turns out, is the right answer. However, the value of  $g$  depends on the particle under consideration. The upshot is that we should include a term in the Hamiltonian of the form

$$H = -g \frac{q}{2m} \mathbf{S} \cdot \mathbf{B} \quad (1.29)$$

### The g-factor

For fundamental particles with spin  $\frac{1}{2}$  — such as the electron — there is a long and interesting history associated to determining the value of  $g$ . For the electron, this was first measured experimentally to be

$$g_e = 2$$

Soon afterwards, Dirac wrote down his famous relativistic equation for the electron. One of its first successes was the theoretical prediction  $g_e = 2$  for any spin  $\frac{1}{2}$  particle. This means, for example, that the neutrinos and quarks also have  $g = 2$ .

This, however, was not the end of the story. With the development of quantum field theory, it was realised that there are corrections to the value  $g_e = 2$ . These can be calculated and take the form of a series expansion, starting with

$$g_e = 2 \left( 1 + \frac{\alpha}{2\pi} + \dots \right) \approx 2.00232$$

where  $\alpha = e^2/4\pi\epsilon_0\hbar c \approx 1/137$  is the dimensionless *fine structure constant* which characterises the strength of the Coulomb force. The most accurate experimental measurement of the electron magnetic moment now yields the result

$$g_e \approx 2.00231930436182 \pm 2.6 \times 10^{-13}$$

Theoretical calculations agree to the first ten significant figures or so. This is the most impressive agreement between theory and experiment in all of science! Beyond that, the value of  $\alpha$  is not known accurately enough to make a comparison. Indeed, now the measurement of the electron magnetic moment is used to *define* the fine structure constant  $\alpha$ .

While all fundamental spin  $\frac{1}{2}$  particles have  $g \approx 2$ , this does not hold for more complicated objects. For example, the proton has

$$g_p \approx 5.588$$

while the neutron — which of course, is a neutral particle, but still carries a magnetic moment — has

$$g_n \approx -3.823$$

where, because the neutron is neutral, the charge  $q = e$  is used in the formula (1.29). These measurements were one of the early hints that the proton and neutron are composite objects.

### 1.5.1 Spin Precession

Consider a constant magnetic field  $\mathbf{B} = (0, 0, B)$ . We would like to understand how this affects the spin of an electron. We'll take  $g_e = 2$ . We write the electric charge of the electron as  $q = -e$  so the Hamiltonian is

$$H = \frac{e\hbar}{2m} \boldsymbol{\sigma} \cdot \mathbf{B}$$

The eigenstates are simply the spin-up  $|\uparrow\rangle$  and spin-down  $|\downarrow\rangle$  states in the  $z$ -direction. They have energies

$$H|\uparrow\rangle = \frac{\hbar\omega_B}{2}|\uparrow\rangle \quad \text{and} \quad H|\downarrow\rangle = -\frac{\hbar\omega_B}{2}|\downarrow\rangle$$

where  $\omega_B = eB/m$  is the cyclotron frequency which appears throughout this chapter.

What happens if we do not sit in an energy eigenstate. A general spin state can be expressed in spherical polar coordinates as

$$|\psi(\theta, \phi)\rangle = \cos(\theta/2)|\uparrow\rangle + e^{i\phi} \sin(\theta/2)|\downarrow\rangle$$

As a check, note that  $|\psi(\theta = \pi/2, \phi)\rangle$  is an eigenstate of  $\sigma^x$  when  $\phi = 0, \pi$  and an eigenstate of  $\sigma^y$  when  $\phi = \pi/2, 3\pi/2$  as it should be. The evolution of this state is determined by the time-dependent Schrödinger equation

$$i\hbar \frac{\partial |\psi\rangle}{\partial t} = H|\psi\rangle$$

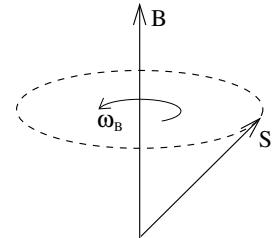


Figure 12:

which is easily solved to give

$$|\psi(\theta, \phi; t)\rangle = e^{-i\omega_B t/2} \left[ \cos(\theta/2) |\uparrow\rangle + e^{i(\phi+\omega_B t)} \sin(\theta/2) |\downarrow\rangle \right]$$

We see that the effect of the magnetic field is to cause the spin to precess about the **B** axis, as shown in the figure.

### 1.5.2 A First Look at the Zeeman Effect

The Zeeman effect describes the splitting of atomic energy levels in the presence of a magnetic field. Consider, for example, the hydrogen atom with Hamiltonian

$$H = -\frac{\hbar^2}{2m} \nabla^2 - \frac{1}{4\pi\epsilon_0} \frac{e^2}{r}$$

The energy levels are given by

$$E_n = -\frac{\alpha^2 mc^2}{2} \frac{1}{n^2} \quad n \in \mathbf{Z}$$

where  $\alpha$  is the fine structure constant. Each energy level has a degeneracy of states. These are labelled by the angular momentum  $l = 0, 1, \dots, n-1$  and the  $z$ -component of angular momentum  $m_l = -l, \dots, +l$ . Furthermore, each electron carries one of two spin states labelled by  $m_s = \pm \frac{1}{2}$ . This results in a degeneracy given by

$$\text{Degeneracy} = 2 \sum_{l=0}^{n-1} (2l+1) = 2n^2$$

Now we add a magnetic field  $\mathbf{B} = (0, 0, B)$ . As we have seen, this results in perturbation to the Hamiltonian which, to leading order in  $B$ , is given by

$$\Delta H = \frac{e}{2m} (\mathbf{L} + g_e \mathbf{S}) \cdot \mathbf{B}$$

In the presence of such a magnetic field, the degeneracy of the states is split. The energy levels now depend on the quantum numbers  $n$ ,  $m_l$  and  $m_s$  and are given by

$$E_{n,m,s} = E_n + \frac{e}{2m} (m_l + 2m_s) B$$

The Zeeman effect is developed further in the [Lectures on Topics in Quantum Mechanics](#).