Funky Electromagnetic Concepts
The Anti-Textbook*
A Work in Progress. See physics.ucsd.edu/~emichels for the latest versions of the Funky Series.
Please send me comments.

Eric L. Michelsen

“But Mr. Faraday, of what use is all this?” - unknown woman
“Madam, of what use is a newborn baby?” - Michael Faraday
“With electromagnetism, as with babies, it’s all a matter of potential.”
- Bill Nye, the Science Guy

* Physical, conceptual, geometric, and pictorial physics that didn’t fit in your textbook.
Instead of distributing this document, please link to physics.ucsd.edu/~emichels/FunkyElectromagneticConcepts.pdf.
Please cite as: Michelsen, Eric L., Funky Electromagnetic Concepts, physics.ucsd.edu/~emichels, 7/7/2017.
“Remarkably clear and easy to reproduce.” -- physics graduate student.
“Finally, a systematic approach to boundary value problems.” -- physics graduate student.


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<td>Bohr magneton</td>
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**Other values:**

Jansky (Jy), flux and spectral density $10^{-26} \text{ W m}^2/\text{Hz} = 10^{-23} \text{ erg/s cm}^2/\text{Hz}$
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1 Introduction

Why Funky?

The purpose of the “Funky” series of documents is to help develop an accurate physical, conceptual, geometric, and pictorial understanding of important physics topics. We focus on areas that don’t seem to be covered well in most texts. The Funky series attempts to clarify those neglected concepts, and others that seem likely to be challenging and unexpected (funky?). The Funky documents are intended for serious students of physics; they are not “popularizations” or oversimplifications.

Physics includes math, and we’re not shy about it, but we also don’t hide behind it.

Without a conceptual understanding, math is gibberish.

This work is one of several aimed at graduate and advanced-undergraduate physics students. Go to http://physics.ucsd.edu/~emichels for the latest versions of the Funky Series, and for contact information. We’re looking for feedback, so please let us know what you think.

How to Use This Document

This work is not a textbook.

There are plenty of those, and they cover most of the topics quite well. This work is meant to be used with a standard text, to help emphasize those things that are most confusing for new students. When standard presentations don’t make sense, come here.

If you don’t understand something, read it again once, then keep reading.

Don’t get stuck on one thing. Often, the following discussion will clarify things.

You should read all of this introduction to familiarize yourself with the notation and contents. After that, this work is meant to be read in the order that most suits you. Each section stands largely alone, though the sections are ordered logically. Simpler material generally appears before more advanced topics. You may read it from beginning to end, or skip around to whatever topic is most interesting.

The index is not yet developed, so go to the web page on the front cover, and text-search in this document.

What’s Wrong With Existing Electromagnetic Expositions?

They’re not precise enough with their definitions. Usually, when there appears to be an obvious contradiction, it is a confusion of definitions. Many widely used references have terribly unclear definitions, and one purpose of these notes is to help resolve them. Also, many texts are not visual or graphical enough. They rely way too much on algebra or advanced math, and not enough on insight.

My Story

The Funky series of notes is the result of my going to graduate school in physics after 20 years out of school. Although I had been an engineer all that time, most of my work involved software and design architectures that are far removed from fundamental science and mathematics. I expected to be a little rusty, but I found that the rust ran deeper than I realized.

There are many things I wish I had understood better while taking my classes (first at San Diego State University, then getting my PhD at University of California, San Diego). The Funky series is my attempt to help other students acquire a deeper understanding of physics.

Thank You

I owe a big thank you to many professors at both SDSU and UCSD, for their generosity, even when I wasn’t a real student: Dr. Herbert Shore, Dr. Peter Salamon, Dr. Arlette Baljon, Dr. Andrew Cooksy, Dr. George Fuller, Dr. Tom O’Neil, Dr. Terry Hwa, and others.
Notation

[Square brackets] in text indicates asides that can be skipped without loss of continuity. They are included to help make connections with other areas of physics.

\[ \text{arg } A \] for a complex number \( A \), \( \text{arg } A \) is the angle of \( A \) in the complex plane; i.e., \( A = |A|e^{i\text{arg } A} \).

[Interesting points that you may skip are “asides,” shown in square brackets, or smaller font and narrowed margins. Notes to myself may also be included as asides.]

Common misconceptions are sometimes written in dark red dashed-line boxes.

Mnemonics (memory aids) or other tips are given in green boxes.
2 Circuits

Circuits Reference Desk

Which end of a resistor is positive? An inductor? A capacitor? A diode? A battery? It all comes down to two simple conventions: the passive convention, and the active convention. We start with the resistor, then proceed to the more complicated devices.

These principles extend directly to AC analysis, using phasors and complex impedance.

\[ v = iR \]

\[ v = L \frac{di}{dt} \]

\[ v = C \frac{dv}{dt} \]

\[ i = I_0 \left( e^{\frac{qV}{kT}} - 1 \right) \]

\[ v = \text{constant} \]

\[ i = \text{arbitrary} \]

**Figure 2.1** Reference polarities for all components, except sources, follow the passive convention. The reference polarity is reversed for the battery (a source) compared to all other components; it uses the active convention.

For any circuit element, conventions define a reference polarity for the voltage, and also a reference direction for the current.

We must write our I-V equations consistently with those choices.

**For a resistor:** the current always flows from + to −, so we choose reference directions consistent with that (diagram above, left). This allows us to write Ohm’s law without minus signs: \( v = iR \). All other passive devices follow this same convention. Therefore, when the power \( P = vi \) is positive, the device is consuming power from the circuit.

Sources, both voltage and current sources, always use the active convention: reference current flows out of the reference positive voltage. Therefore, when the power \( P = vi \) is positive, the source is supplying power to the circuit.

For an arbitrary circuit, we may not know ahead of time which end of a resistor will end up being +. For example:

**Figure 2.2** Two valid choices of reference directions for the same circuit.

We don’t know, without being given numbers and doing the circuit analysis, whether the middle resistor’s current will flow up or down. No problem: we just choose an arbitrary polarity (this defines both voltage and current, since their relationship is fixed by the conventions: reference current flows from reference + to reference −). We do the circuit analysis assuming this polarity. Note that:

Ohm’s law applies for both positive and negative voltages and currents.

If we find that the voltage (and therefore current) is negative, it just means that the current is really flowing opposite to our reference choice. Note, however, that resistors always consume power from the circuit:
For a capacitor: Things are a tad messier, because current does not always flow from + to –. A capacitor stores energy in its electric field. If we increase the voltage on the capacitor from zero, the capacitor is drawing energy from the rest of the circuit (it is charging). In this case, it is qualitatively similar to a resistor, and its current flows from + to –. Therefore, to be consistent with resistors, we define our reference directions for this case: reference current flows from reference + to reference – (just like a resistor). But wait! The current through a capacitor is not related to the polarity of voltage across it; the current is related to the rate of change of voltage:

\[ i = C \frac{dv}{dt} \quad \text{units: } \frac{C}{s} = \frac{(V)}{(V/s)} \]

If the voltage is negative, but increasing (becoming less negative), the current through the capacitor is still positive. Thus the \( I-V \) equation above is always valid: charging or discharging. When a capacitor is discharging, either its voltage is + and its current is –, or its voltage is – and its current is +. Either way, the capacitor is delivering energy to the circuit, and temporarily acts more like a battery than a resistor. But we can’t change our reference directions on the circuit based on the charging/discharging state of the capacitor from moment to moment. Also, either way, the capacitor’s power “consumed” is

\[ P = VI \quad \text{where negative power means it supplies energy to the circuit.} \]

For an inductor: things are similar, but we exchange “voltage” and “current”, and replace “electric” with “magnetic.” Again, the current does not always flow from + to –. An inductor stores energy in its magnetic field. If we increase the current from zero, the inductor is drawing energy from the rest of the circuit (loosely, it is “charging”). In this case, it is qualitatively similar to a resistor, and its current flows from + to –. Again, to be consistent with resistors, we define our reference directions for this case: reference current flows from reference + to reference – (just like a resistor). But wait! The voltage across an inductor is not related to direction of the current through it; the voltage is related to the rate of change of the current. Therefore, if the current is negative, but increasing (becoming less negative), the voltage across the inductor is still positive. This allows us to write a single \( I-V \) equation for both cases:

\[ v = L \frac{di}{dt} \quad \text{units: } V = \frac{\text{flux-linkages}}{\text{ampere}} \left( \frac{\text{ampere}}{s} \right) \]

When an inductor is “discharging,” the current is decreasing, and the inductor is supplying energy to the circuit. Now it is acting more like a battery than a resistor. But we made our reference choice for the inductor, and we must stick with it. The above \( I-V \) equation is valid at all times. Again,

\[ P = VI \quad \text{where negative power means it supplies energy to the circuit.} \]

Note that when the inductor current is increasing (either becoming more positive or less negative), \( v \) is positive. When the current is decreasing (either becoming less positive or more negative), \( v \) is negative, which means the + reference terminal is really at negative voltage with respect to the – terminal. In all cases, the above equations are correct. We achieved that consistency by defining a single reference polarity.

For a diode: Resistors, capacitors, and inductors are all “symmetric” or “unpolarized” devices: you can reverse the two leads with no effect. Diodes, in contrast, are polarized: one lead is the “anode”; the other is the “cathode.” You must connect them properly. The reference voltage is defined always with + on the anode (and therefore – on the cathode); reference current flows from + to –. Diodes always consume power (like resistors do). Consistently with that, these conventions require that

\[ P = VI \quad \text{is always positive: the diode consumes power.} \]

For a battery: Batteries usually supply energy to the circuit, so we define them as having positive power when they do so (the opposite of all other devices here). This requires the opposite reference directions:

\[ \text{Batteries use the opposite convention from other devices:} \]

\[ \text{reference current flows through the battery from reference – to reference +.} \]
However, it is possible to force current “backwards” through a battery, and then it will consume energy from the circuit (as demanded by the fundamental definitions of voltage and current). This is how we recharge a rechargeable battery. Thus:

\[ P = VI \] where positive power \( \Rightarrow \) battery supplies energy to the circuit; negative power \( \Rightarrow \) battery consumes energy from the circuit.

For a transformer: Transformers are a new breed, because they have 4 terminals, rather than two:

\[ v_1 = M \frac{di_2}{dt} \quad \text{and} \quad v_2 = M \frac{di_1}{dt} \]
\[ v_2 = Nv_1 \quad \text{and} \quad i_2 = \frac{1}{N} i_1 \]

An ideal transformer relates the voltages and currents on both sides.

The dots on the two windings define the reference directions and polarities, as shown in the diagram above. Essentially, the reference directions are the same as for an inductor, except that \( v_1 \) depends on \( i_2 \), and \( v_2 \) depends on \( i_1 \). This implies that when \( v_1 \) is positive on the dot-side, \( v_2 \) is also positive on the dot-side.

In an ideal transformer, all the magnetic flux, \( \Phi \), passes through every turn of both the primary and secondary windings. From Faraday’s law, \( V = -\frac{d\Phi}{dt} \), applied to each turn of wire:

\[ v_1 = -N_1 \frac{d\Phi}{dt}, \quad v_2 = -N_2 \frac{d\Phi}{dt} \Rightarrow v_2 = \frac{N_2}{N_1} v_1 = Nv_1 \quad \text{where} \quad N = \frac{N_2}{N_1} \]

Also, an ideal transformer has a highly magnetizeable core such that the flux, \( \Phi \), requires virtually no current to create it. This means that the primary and secondary currents must cancel each other, leaving nearly zero MMF (magneto-motive force). Therefore:

\[ N_1 i_1 = N_2 i_2 \quad \text{or} \quad i_2 = \frac{1}{N} i_1 \quad \text{where} \quad N = \frac{N_2}{N_1} \quad \text{as before} \]

The secondary voltage varies as the turns ratio, \( N \).
The secondary current varies as the inverse of the turns ratio.

This implies that the power into the primary equals the power out of the secondary. Unlike an inductor, an ideal transformer does not store energy. Instead, a transformer transfers energy from one side of the transformer to the other. The direction of energy transfer depends on the circuits to which the two transformer sides are connected. Details are beyond the scope of this section.

**Brief Note on Phasor Analysis**

This section assumes you are familiar with phasor analysis, also known as Fourier mode analysis. The idea is to consider one pure sinusoidal frequency at a time. Since the circuit response is linear with the excitation, a sum of sinusoidal excitations results in a response equal to the sum of the individual sinusoidal responses.

For capacitors, we have (using \( \exp (+i\omega t) \) time dependence, which is standard for circuits):

\[ i(t) = C \frac{dv}{dt} \Rightarrow \tilde{i} = i\omega\tilde{v} \quad \text{where} \quad \tilde{i}, \tilde{v} \text{ are the phasors for current and voltage} \]

The phase of the derivative is positive, which means the derivative leads the original function. In this case, the current leads the voltage, or equivalently, the voltage lags the current (below left).
(Left) In a capacitor, voltage lags current. (Right) In an inductor, current lags voltage.

For inductors, things are reversed:

\[ v(t) = L \frac{di}{dt} \Rightarrow \vec{v} = i \omega L \vec{i} \]

where \( \vec{i}, \vec{v} \) are the phasors for current and voltage.

The voltage leads the current, or the current lags the voltage (above right).

You can remember the lead/lag relationships for circuit elements mnemonically as follows:

- Capacitors oppose a change in voltage, so the voltage lags the current.
- Inductors oppose a change in current, so the current lags the voltage.

Similarly:

You can remember the lead/lag relationships for derivatives as follows:

- Derivatives show where you’re going before you get there, so they lead the function.
- Integrals sum up where you’ve been, and so lag the function.
### 3 Classical Electromagnetics

#### Just For Reference: Faraday’s Law

What is the meaning of the minus sign in Faraday’s law? Faraday’s law relates the induced voltage in a circuit surrounding a magnetic field to the rate of change of that field:

\[ V = -\frac{d\Phi}{dt} \]

The minus sign can only be interpreted with respect to a standard set of reference directions and polarities. As with reference directions in electric circuits (see Circuits Reference Desk elsewhere), a **reference direction** (of a current, B-field, or E-field) is the direction which is called positive in the equations. A **reference polarity** (of a voltage) is the polarity which is called positive in the equations. The actual polarity may be the same as our reference choice, or it may be opposite. We may not know until we solve some equations. But there is no problem either way, because if the actual polarity is opposite our reference choice, it will simply have a negative value.

![Reference directions for flux Φ, current, and reference polarity for voltage.](image)

**Figure 3.1** Reference directions for flux Φ, current, and reference polarity for voltage.

Now, Faraday’s law: the reference direction for current is that which would produce the reference direction for Φ, and Φ is in the same direction as B. In the diagram above, B and Φ are positive in the z-direction (out of the page). The reference direction for the current is counter-clockwise, which produces (by the right hand rule) a B-field out of the page. We can lump the resistance of the loop into a single equivalent resistor. The reference polarity for the voltage must be consistent with the reference direction for the current, and this forces the choice shown. We now have a consistent set of reference directions and polarity for all four of the flux, B-field, current, and voltage. Positive voltage makes positive B-field.

This defines the meaning of the minus sign in Faraday’s law. Lenz’ law says that if the B-field changes, the voltage induced will try to drive a current that produces a B-field which opposes the change. This is the minus sign: decreasing B (negative \(d\Phi/dt\)) causes positive voltage, which boosts B. Increasing B (positive \(d\Phi/dt\)) causes negative voltage, which reduces B.

The minus sign emphasizes that if the induced voltage reinforced the change in B-field, that would induce more voltage, further changing the B-field, which induces more voltage, further changing the B-field, in a never ending death spiral of infinite current and B-field.

### Stunning Phasors and Fourier Space

Phasors are a convenient way to mathematically represent oscillations. Phasors and Fourier space are used heavily in EM waves and propagation, as well as classical mechanics, quantum mechanics, and any other physics that involves oscillations. Phasors are also used extensively in engineering. Virtually all physics curricula use phasors, though many do not explain the concept, or use the word “phasor.” As a result, many physics students are limited by not clearly understanding phasors. A full understanding of phasors includes understanding ratios of phasors (e.g., impedance), and simple extensions of other concepts, such as a complex propagation vector, complex permittivity, and complex index of refraction. Phasors also provide a simple but sturdy foundation on which to build more advanced concepts: e.g., a Fourier transform is a phasor-valued function of frequency; a quantum wave-function can be considered a phasor-valued function of space.
Working with phasors is also called working in Fourier space or in Fourier modes. A phasor can also be called a complex amplitude [Gri E&M p???]. (Quantum mechanics uses the term “complex amplitude” for a complex number whose relationship to a sinusoid is somewhat abstract.)

This section requires that you understand complex numbers in both polar and rectangular form, and the basic calculus of complex functions of real variables.

We start by noting that any real-valued sinusoid is fully characterized by 3 numbers: amplitude, phase, and frequency:

\[ c(t) = C \cos(\omega t + \theta), \quad C \text{ real} \]

**A phasor** is a complex number that characterizes the amplitude and phase of a sinusoid.

A phasor says nothing about its frequency. You must know the frequency from some other condition. Combining two phasors by simple addition only makes sense if they refer to sinusoids of the same frequency. However, phasors of different frequencies are often combined by inserting the time dependence explicitly before combining (quantum mechanics does this routinely).

A phasor \( A \) (a complex number) corresponds to the sinusoid (in the engineering time convention):

\[ a(t) = |A| \cos(\omega t + \text{arg} A) = \text{Re}\{A e^{i\omega t}\}, \quad \text{where arg } A = \text{the complex angle of } A. \]

**The sign of the times:** For engineers (including electromagnetics) and classical physics [M&T, Tay], and for AC circuit analysis, the time dependence is \( e^{i\omega t} \). For most physicists in quantum mechanics and electromagnetics, the time dependence is \( e^{-i\omega t} \). You can remember this mnemonically by thinking that for wave physicists, time goes backwards. The polarity of the exponent is purely conventional, and has no physical significance. This work shows pictures for both time conventions.

The magnitude of the phasor is exactly the (real) amplitude of the sinusoid, and the complex angle of the phasor is exactly the phase of the sinusoid, i.e. the angle of the cosine at \( t = 0 \). The geometric interpretation of a phasor is that of a rotating stick, which casts a shadow on the horizontal axis. The length of the shadow is a sinusoidal function of time, with amplitude \( |A| \), and starting at an angle \( \text{arg} A \):

### Figure 3.2

For engineers, the stick rotates counter-clockwise, as shown, per \( e^{i\omega t} \). For wave physicists, the stick usually rotates clockwise (opposite to that shown), per \( e^{-i\omega t} \).

We can also view the rotation at frequency \( \omega \) as complex multiplication by \( e^{i\omega t} \) (engineering), or \( e^{-i\omega t} \) (physics). Recall that multiplication by a unit-magnitude complex number simply rotates another complex number in the complex plane. Now imagine the unit-magnitude angle is not fixed, but changes linearly with time, i.e. multiply not by a fixed complex angle \( \theta \), but by an increasing angle \( \omega t \). \( \omega \) is the angular frequency, in rad/s. When we multiply some complex number \( r e^{i\theta} \) by \( e^{i\omega t} \), we get a complex function of time that rotates continuously around the origin in the complex plane. The magnitude of the result is fixed, because \( |e^{i\omega t}| = 1 \) at all times. But the angle of the result increases with time, at the rate \( \omega \).
The angular frequency is not constrained to be positive; it can just as well be negative. Rotation by a negative frequency rotates in the clock-wise direction, rather than counter clockwise. Hence, both positive and negative frequencies occur in complex rotations, and in signal analysis.

Recall that any linear combination of sinusoids, of arbitrary amplitude and phase (but identical frequency, of course), is another sinusoid (at that frequency). The beauty of phasors is that the phasor representing the sum of two sinusoids is simply the (complex) sum of the phasors representing the original sinusoids (addends). The graphical demonstration of this is both simple, and illuminating:

Let phasor \( A + B \). Then \( a(t) = \text{Re} \{ A e^{-i\omega t} + B e^{-i\omega t} \} = \text{Re} \{ (A_r + i A_i) e^{-i\omega t} \} = A_r \cos \omega t + A_i \sin \omega t \).

Thus we see that the real and imaginary parts of the phasor \( A \) are exactly the in-phase and quadrature components of the sine wave.

We use complex numbers to represent sinusoids because the arithmetic (and some calculus) of complex numbers (2D vectors) is the same as the arithmetic of adding sinusoids.

**Direction of the wave-vector:** For traveling waves, we can see that the wave-vector points in the direction of propagation, by considering a point of constant phase on the wave. For constant phase, we must have

\[
k \cdot x - \omega t = \text{const}
\]

\( \Rightarrow \) \( k \) points in direction of propagation.
because as \( t \) increases, so must \( \mathbf{k} \cdot \mathbf{x} \), and therefore the point of constant phase \((\mathbf{x} \rightarrow \mathbf{x} + d\mathbf{x})\) must move in the same direction as \( \mathbf{k} \) points.

Even in the engineering convention, with time evolution given by \( \exp(+i\omega t) \), the wave-vector \( \mathbf{k} \) still points in the direction of propagation. The condition for propagation is \( \omega t - \mathbf{k} \cdot \mathbf{x} = \text{const} \), which can be brought into the physics-convention form by absorbing a minus sign into the constant.

**Phasor Calculus**

We can easily see that phasors convert differential equations to algebraic equations. This is expected, because phasors are a method of Fourier analysis, which is well known for converting differential to algebraic equations. Let’s take the first and second derivatives of a cosine, in both the old-fashioned real-valued way, and this new-fangled phasor way:

We can take time derivatives of phasors by noting that the time derivative of the real part of a complex function equals the real part of the time derivative:

\[
\frac{d}{dt} \text{Re}\{z(t)\} = \text{Re}\left\{ \frac{d}{dt} z(t) \right\}
\]

because

\[
\text{Re}\left\{ \frac{d}{dt} z(t) \right\} = \text{Re}\left\{ \frac{d}{dt} z_r(t) + i \frac{d}{dt} z_i(t) \right\} = \frac{d}{dt} z_r(t) = \frac{d}{dt} \text{Re}\{z(t)\}.
\]

Then

\[
\frac{d}{dt} \text{Re}\{Ae^{-i\omega t}\} = \text{Re}\left\{ \frac{d}{dt} Ae^{-i\omega t} \right\} = \text{Re}\{ -i\omega Ae^{-i\omega t} \}
\]

In phasor notation:

\[
\frac{d}{dt} A = -i\omega A
\]

For example (in the engineering convention):

\[
\cos(\omega t) \quad A = 1 \quad \leftrightarrow \quad \text{Re}\{Ae^{i\omega t}\} = \text{Re}\{1(\cos \omega t + i \sin \omega t)\} = \cos \omega t
\]

\[
\frac{d}{dt} \cos \omega t = -\omega \sin \omega t
\]

\[
\frac{d}{dt} A = i\omega A \quad \leftrightarrow \quad \text{Re}\{i\omega Ae^{i\omega t}\} = \text{Re}\{i\omega(\cos \omega t + i \sin \omega t)\} = \text{Re}\{i\omega \cos \omega t - \omega \sin \omega t\} = -\omega \sin \omega t
\]

\[
\frac{d^2}{dt^2} \cos \omega t = -\omega^2 \cos \omega t
\]

\[
\frac{d^2}{dt^2} A = (i\omega)^2 A = -\omega^2 A \quad \leftrightarrow \quad \text{Re}\{-\omega^2 Ae^{i\omega t}\} = \text{Re}\{-\omega^2 (\cos \omega t + i \sin \omega t)\} = -\omega^2 \cos \omega t
\]

We can also have phasors defining both space and time sinusoidal variations. These can be used for traveling waves.

Then the phasor carries the amplitude and phase of the traveling wave, but not its wave-vector \( \mathbf{k} \) (spatial frequency) or temporal frequency \( \omega \). We must be given \( \mathbf{k} \) and \( \omega \) separately. In 1-dimension, the spatial derivative works like this:
Let \( z(t,x) \equiv z_r(t,x) + iz_j(t,x) \).

Then \( \frac{d}{dx} \text{Re}(z(t,x)) = \text{Re}\left\{ \frac{d}{dx} z(t,x) \right\} \) similar to above with \( \frac{d}{dt} \),
and
\[
\frac{d}{dx} \text{Re}(A e^{ikx-i\omega t}) = \text{Re}\left\{ \frac{d}{dx} A e^{ikx-i\omega t} \right\} = \text{Re}\left\{ ikA e^{ikx-i\omega t} \right\}
\]

In phasor notation:
\[
\frac{d}{dx} A = ikA.
\]

In higher dimension space, we replace \( \frac{\partial}{\partial x} \) with \( \nabla \):

Let \( z(t,r) \equiv z_r(t,r) + iz_j(t,r) \).

Then \( \nabla \text{Re}(z(t,r)) = \text{Re}\left\{ \nabla z(t,r) \right\} \) similar to above with \( \frac{d}{dx} \).

Then \( \nabla \text{Re}(A e^{ikr-i\omega t}) = \text{Re}\left\{ \nabla A e^{ikr-i\omega t} \right\} = \text{Re}\left\{ ikke^{ikr-i\omega t} \right\} \)

In phasor notation:
\[ \nabla A = i k A. \]

**Time Averages**

It is often useful to compute the time average of the product of 2 sinusoids. E.g., the time average of the Poynting vector gives the effective radiation power density in W/m\(^2\) (or power per unit area). The time average of two sinusoids does not depend on the absolute phase of either; it depends only on the relative phase of the two. Specifically,

\[
\langle a(t)b(t) \rangle_t \equiv \frac{1}{2} \text{Re}\{AB^*\}.
\]

One way to see this is to decompose the phasors into the cosine (real) and sine (imaginary) components.

\[
a(t) = A_r \cos \omega t + A_i \sin \omega t \quad \quad b(t) = B_r \cos \omega t + B_i \sin \omega t
\]

\[
\langle a(t)b(t) \rangle_t = \left\langle \left( A_r \cos \omega t + A_i \sin \omega t \right) \left( B_r \cos \omega t + B_i \sin \omega t \right) \right\rangle_t
\]

\[
= \left\langle A_r B_r \cos^2 \omega t + A_r B_i \cos \omega t \sin \omega t + A_i B_r \cos \omega t \sin \omega t + A_i B_i \sin^2 \omega t \right\rangle_t
\]

\[
= \left\langle A_r B_r \cos^2 \omega t \right\rangle_t + \left\langle A_r B_i \cos \omega t \sin \omega t \right\rangle_t + \left\langle A_i B_r \cos \omega t \sin \omega t \right\rangle_t + \left\langle A_i B_i \sin^2 \omega t \right\rangle_t
\]

Only the cos-cos and sin-sin terms contribute to the time average, because the time average of cos-sin is zero (they are orthogonal functions; sin is odd, cos is even). Therefore,

\[
\langle \cos^2 \omega t \rangle_t = \langle \sin^2 \omega t \rangle_t = \frac{1}{2}
\]

\[
\langle a(t)b(t) \rangle_t = \left\langle A_r B_r \cos^2 \omega t \right\rangle_t + \left\langle A_r B_i \sin^2 \omega t \right\rangle_t = \frac{1}{2} A_r B_r + \frac{1}{2} A_i B_i = \frac{1}{2} \text{Re}\{AB^*\}
\]

[Notice that \( \text{Re}\{AB^*\} \) is analogous to the dot-product of two spatial vectors: it is the component of \( B \) parallel to \( A \), times \( |A| \), i.e. \( AB \cos \theta \), where \( \theta = \text{arg } B - \text{arg } A \). Only the parallel components contribute to the time average.]
Figure 3.5 Phasors showing average power: ??

If you’d rather grind through the integration, we can demonstrate the time average formula that way. We start with the sinusoids as real-valued cosines, and later switch back to the complex formula (after the integration):

$$a(t) = |A| \cos(\omega t + \arg A) = \Re\{Ae^{-i\omega t + \arg A}\}$$

$$b(t) = |B| \cos(\omega t + \arg B) = \Re\{Be^{-i\omega t + \arg B}\}$$

$$\langle a(t)b(t) \rangle_t = \int_{\text{period}} dt |A| \cos(\omega t + \arg A)|B| \cos(\omega t + \arg B)$$

$$= |AB| \int_{\text{period}} dt \cos(\omega t + \arg A) \cos(\omega t + \arg B)$$

$$= |AB| \int_{\text{period}} dt \cos(\omega t + (\arg A - \arg B)) \cos(\omega t)$$

The last equation is because the integral is over a full period, so shifting the starting and ending point by a fixed time interval, or angle, doesn’t change its value. This demonstrates that the time average depends only on the phase difference between A and B, and not their absolute phases. Finally,

Use $\cos(c + d) = \cos c \cos d - \sin c \sin d$ where $c = \omega t$, $d = (\arg A - \arg B)$

$$\langle a(t)b(t) \rangle_t = |AB| \int_{\text{period}} dt \cos \omega t \cos(\arg A - \arg B) - \sin \omega t \sin(\arg A - \arg B) \cos(\omega t)$$

$$= \left[ |AB| \cos(\arg A - \arg B) \int_{\text{period}} dt \cos^2(\omega t) \right]$$

$$- \left[ |AB| \sin(\arg A - \arg B) \int_{\text{period}} dt \sin \omega t \cos(\omega t) \right]$$

$$= \frac{1}{2} |AB| \cos(\arg A - \arg B) = \frac{1}{2} \Re \left[ |AB| e^{i(\arg A - \arg B)} \right] = \frac{1}{2} \Re \left[ |A| e^{i\arg A} |B| e^{-i\arg B} \right]$$

$$= \frac{1}{2} \Re \{AB^*\} \quad QED$$

Polarization Vector

This section assumes you understand phasors, and the decomposition of a sinusoid into “in-phase” and “quadrature” parts (see “Phasors,” above). We start with an overview, and then fill in the details.

Overview: A polarization vector in general applies to any vector (or other multi-component object) which travels in space, and varies sinusoidally in time and space. It gives the relative amplitude and phase
of each component of the vector. By convention for EM waves, the polarization vector, ε, gives the relative amplitudes and phases of the E-field components (E_x, E_y, E_z) for the wave.

In applications where the A field is most relevant, we write a polarization vector for the A-field, instead. In this case, ε is gauge–dependent. In some gauges, it has unexpected components, including a “longitudinal” component, which points along the propagation direction. This is a gauge artifact, and the polarization vectors for E and B-fields (in vacuum) always have no component along the propagation direction.

In principle, one could write a polarization vector for a B-field, but it’s not usually done since it is just a right-hand rotation of the E-field polarization vector about the propagation direction by 90°.

[Note that for gravity waves, where the propagating field is the 4 × 4 metric tensor perturbation field, h_μν, the polarization “object” is a 4 × 4 polarization tensor.]

For a propagating EM wave, the polarization vector ε gives two things:

1. the E-field direction (polarization) of the EM wave at any time
2. the fraction of the total EM wave intensity (power density) carried by each component of polarization

The polarization vector ε is a set of 3 phasors, that describe the sinusoidal oscillations of each component (E_x, E_y, E_z) of the E-field. So we can immediately write the form of the E-field at a point as a function of time (in several different notations):

\[
\mathbf{E} \propto \text{Re}\left(\mathbf{e} e^{-i\omega t}\right) = \text{Re}\left(\begin{bmatrix} E_x e^{-i\omega t} \\ E_y e^{-i\omega t} \\ E_z e^{-i\omega t} \end{bmatrix}\right) = \begin{bmatrix} \text{Re}\{E_x e^{-i\omega t}\} \\ \text{Re}\{E_y e^{-i\omega t}\} \\ \text{Re}\{E_z e^{-i\omega t}\} \end{bmatrix}. \]

It is possible to represent any polarization (linear or elliptical) in any propagation direction with a 3D complex vector. Since EM waves are transverse, the E-field, and thus the polarization vector, are perpendicular to the wave-vector \( \mathbf{k} \), i.e., \( \mathbf{e} \cdot \mathbf{k} = 0 \):

![Figure 3.6](image)

**Figure 3.6** (Left) Wave vector \( \mathbf{k} \) in an arbitrary direction, and some possible real polarization vectors. (Right) Wave vector \( \mathbf{k} \), and the real (blue) and imaginary (red) parts of a complex polarization vector.

To keep the x, y, and z intensities normalized with respect to total intensity, the polarization vector is usually normalized to unit magnitude:

\[
\mathbf{e} = \left(\begin{bmatrix} E_x \\ E_y \\ E_z \end{bmatrix}\right) = \mathbf{e}_x \hat{x} + \mathbf{e}_y \hat{y} + \mathbf{e}_z \hat{z}, \quad \text{where} \quad E_x, E_y, E_z \text{ are phasors (complex numbers), and}
\]

\[
|E_x|^2 + |E_y|^2 + |E_z|^2 = 1.
\]
Note, however, there is a subtle distinction between the amplitudes of spatially perpendicular components, and the instantaneous maximum magnitude of the electric field, $E_{\text{max}}$. For a linearly polarized wave, say halfway between the $x$ and $y$-directions, $E_{\text{max}} = (\sqrt{2})E_x$, and the time averaged intensity $\langle I \rangle \sim E_{\text{max}}^2/2$. However, for the same amplitude of perpendicular electric field components, but phase shifted to make circular polarization, $E_{\text{max}} = E_x = E_y$, and $\langle I \rangle \sim E_{\text{max}}^2$. Note that the intensity is the same for both waves: perpendicular component intensities always add; it is $E_{\text{max}}$ which is different.

For linearly polarized waves, $\langle I \rangle \sim E_{\text{max}}^2/2$, whereas for circular waves, $\langle I \rangle \sim E_{\text{max}}^2$.

The real part of the polarization vector is the “in-phase” $E$-field parts. The imaginary part is the time-quadrature $E$-field parts, i.e. is $90^\circ$ out of phase with the real part. Any polarization vector for a propagating wave in a given direction $\mathbf{k}$ can be written as a linear combination of two basis polarization vectors for that $\mathbf{k}$.

**Figure 3.7** EM wave propagating out of the page: sample polarizations

Non-plane waves can be thought of as an infinite set of tiny plane wave fronts propagating in different directions; therefore, non-plane waves have a separate polarization vector at each point. I.e., the polarization vector is a function of space, $\mathbf{\varepsilon}(\mathbf{r})$. Since the reference time for the sinusoidal phase is arbitrary, we conventionally take the first non-zero component of $\mathbf{\varepsilon}$ to be real, which sets the reference time, and define the subsequent components relative to that.

**Details:** We restrict our attention to a single point in space, through which an EM wave is propagating. Since we are at an infinitesimal region of space, any shape wave front may be viewed as a plane wave. For a simple, plane-polarized wave, the time-average intensity $P$ (W/m$^2$) and E-field amplitude $E$ are related by *(the $\varepsilon$ below are permittivity, not polarization vector):*

\[
\begin{align*}
\text{(SI)} & \quad P = \frac{1}{2} \sqrt{\frac{\varepsilon}{\mu}} E^2 \quad \Rightarrow \quad E = \sqrt{2 \frac{\mu}{\varepsilon} P} \quad \text{[Jac 7.13 p298]} \\
\text{(gaussian)} & \quad P = \frac{c}{4\pi} E^2 \quad \Rightarrow \quad E = \frac{4\pi}{c} \sqrt{P} \quad \text{[L&L p120]} \\
\text{In general:} & \quad E = m \sqrt{P}, \quad m = \sqrt{2 \frac{\mu}{\varepsilon}} \quad \text{(SI)} \quad \text{or} \quad m = \frac{4\pi}{c} \quad \text{(gaussian)}
\end{align*}
\]

Consider an EM wave traveling in an arbitrary direction, and with arbitrary polarization. At any given time, the wave has an $E$-field with components in the $x$, $y$, and $z$ directions:

\[
\mathbf{E}(t) = E_x(t)\hat{x} + E_y(t)\hat{y} + E_z(t)\hat{z}
\]

Each component is a sinusoid of frequency $\omega = c|\mathbf{k}|$, and so may be represented by a phasor (recall that a phasor is a complex number which represents the amplitude and phase of a sinusoid; see *Phasors* above):
The propagation vector is \( \mathbf{k} \) (not to be confused with \( \mathbf{k} \equiv \mathbf{z} \)-hat), whose magnitude is the spatial frequency in rad/m, and whose direction is the propagation direction. For a given E-field component \( j \in \{x, y, z\} \), the intensity (power/area) carried in the \( \mathbf{k} \) direction is proportional to \( E_j^2 \), where \( E_j \) is the (real) amplitude:

\[
P_x = m^2 E_x^2 = |e_x|^2 P_{\text{total}} \quad P_y = m^2 E_y^2 = |e_y|^2 P_{\text{total}} \quad P_z = m^2 E_z^2 = |e_z|^2 P_{\text{total}}
\]

\[
P_{\text{total}} = P_x + P_y + P_z \quad \text{W/m}^2 \quad \Rightarrow \quad |e_x|^2 + |e_y|^2 + |e_z|^2 = 1 \quad \text{[Jac 7.13 p298]}
\]

So \( |e_i|^2 \) is the fraction of the total wave intensity carried by the \( E_i \) component, etc.

It is instructive to consider the real and imaginary parts of \( \varepsilon \) separately:

\[\varepsilon = \varepsilon_R + i\varepsilon_I \quad \text{where} \quad \varepsilon_R \text{ and } \varepsilon_I \text{ are real vectors} \]

\( \varepsilon_R \) is the direction of the E-field at the reference time, \( t = 0 \), and all integer oscillation periods after, \( t = nT = 2\pi n/\omega \) (Figure 3.6, right diagram, blue arrow). \( \varepsilon_I \) is the direction of the E-field at quadrature times, \( t = (n + 1/4)T \) (Figure 3.6, right diagram, red arrow). The E-field is always perpendicular to \( \mathbf{k} \), so both \( \varepsilon_R \) and \( \varepsilon_I \) lie in the plane perpendicular to \( \mathbf{k} \).

\[\varepsilon_R \cdot \mathbf{k} = \varepsilon_I \cdot \mathbf{k} = 0 \quad \Rightarrow \quad \varepsilon \cdot \mathbf{k} = 0 + 0i = 0\]

**Basis polarization vectors:** Since EM waves are transverse, the E-field (and therefore polarization vector) lies in a plane perpendicular to \( \mathbf{k} \). This plane is a 2-D space, and all vectors in that plane (for a given direction of \( \mathbf{k} \)) can be written as the linear combination of 2 basis vectors. For complex vectors like polarization vectors, the basis vectors can be either real or complex. We always choose the basis vectors to be orthonormal, i.e.

\[e_1 \cdot e_1 = e_2 \cdot e_2 = 1 \quad e_1 \cdot e_2 = 0, \quad \text{where} \quad a \cdot b = a_x^* b_x + a_y^* b_y + a_z^* b_z\]

[The dot product of complex vectors is just like a quantum inner product.] The only two bases you ever see are linear polarization bases, and circular polarization bases. To simplify the discussion, we now focus on propagation in the \( z \)-direction (\( \mathbf{k} = k \mathbf{e}_z \)). There are 2 orthogonal directions of linear polarization, \( x \) and \( y \). Therefore, our linear polarization basis vectors are simply:

\[e_1 = (1, 0, 0) \quad e_2 = (0, 1, 0) \quad \text{linear polarization basis vectors} \]

For \( z \) propagation, the plane of the E-field is the \( x-y \) plane; thus every polarization vector for \( z \) propagation can be written as a (possibly complex) combination of \( e_1 \) and \( e_2 \). The polarization vector for 100% \( x \) polarization is just \( \varepsilon = e_1 = (1, 0, 0) \). For 100% \( y \) polarization, \( \varepsilon = e_2 = (0, 1, 0) \). For linear polarization at \( 45^\circ \), \( \varepsilon = (1/\sqrt{2}, 1/\sqrt{2}, 0) \), which means \( 1/2 \) the power is carried in \( E_x \) and half in \( E_y \).

For right-hand-circular polarization (RHC), \( \mathbf{E}(t) \) rotates counter clockwise (right hand rule applied to \( \mathbf{k} \)). This means \( \mathbf{E}(t = 0) \) points in the \( x \) direction, so \( \varepsilon_R \sim (1, 0, 0) \). \( \mathbf{E}(t = T/4) \) points in the \( y \) direction, so \( \varepsilon_I \sim (0, 1, 0) \). Then

\[\varepsilon - \varepsilon_R + i\varepsilon_I = (1, 0, 0) + i(0, 1, 0) = (1, i, 0). \quad \text{Normalizing,} \quad \varepsilon = \frac{1}{\sqrt{2}}(1, i, 0)\]

For LHC, \( \mathbf{E}(t = T/4) \) points in the \( -y \) direction, so \( \varepsilon = (1/\sqrt{2}, -i/\sqrt{2}, 0) \). We have just derived the circular polarization basis vectors:

\[\text{For RHC,} \quad \varepsilon_R = (1, 0, 0) \quad \varepsilon_I = (0, 1, 0) \]

\[\text{For LHC,} \quad \varepsilon_R = (1/\sqrt{2}, -i/\sqrt{2}, 0) \quad \varepsilon_I = (1/\sqrt{2}, i/\sqrt{2}, 0)\]
\[ \mathbf{e}_R = \frac{1}{\sqrt{2}} (1, i, 0) \]
\[ \mathbf{e}_L = \frac{1}{\sqrt{2}} (1, -i, 0) \]

[Beware that in optics, RHC is called “left handed” and LHC is called “right handed.” Beats me.]

**Examples:** Let’s use our polarization basis vectors to find out how much power passes through some polarization filters.

1. Given an elliptical wave with \( \mathbf{e} = (3/5, i4/5, 0) \), what fraction \( f \) of the power will survive a y-polarizing filter? The filter passes only the part of the wave with \( \mathbf{e} = (0, 1, 0) \). Therefore, we “project” our given polarization onto this, i.e., we see what fraction of our wave is \( E_y \)-polarized. This is just like finding how much of a basis vector is in a quantum state. Then we square that, to get intensity:

\[
\begin{align*}
f &= |\mathbf{e}_y \cdot \mathbf{e}|^2 = |(0,1,0) \cdot (3/5,i4/5,0)|^2 = |4/5|^2 = 16/25
\end{align*}
\]

A circular wave, \( \mathbf{e} = (1/\sqrt{2}, i/\sqrt{2}, 0) \), and a 45° linear wave, \( \mathbf{e} = (1/\sqrt{2}, 1/\sqrt{2}, 0) \), would both pass \( 1/2 \) the power.

2. How much of an \( x \) polarized wave survives a chiral filter that passes only RHC?

\[
\begin{align*}
f &= |\mathbf{e}_x \cdot \mathbf{e}|^2 = |(1/\sqrt{2}, i/\sqrt{2}, 0) \cdot (a, b, 0)|^2 = \frac{(a+ib)^2}{\sqrt{2}} = 1/2,
\end{align*}
\]

The answer is the same for any direction of linear polarization (axial symmetry):

\[
\begin{align*}
&= |\mathbf{e}_x \cdot \mathbf{e}|^2 = \left| \frac{1}{\sqrt{2}} \right|^2 = 1/2, \quad \text{since} \quad |\mathbf{e}|^2 = 1 = |a+ib|^2
\end{align*}
\]

Notice that in these cases, linear polarized light went in, but circular polarized light came out.

3. How much of an LHC wave would pass the RHC filter?

\[
\begin{align*}
f &= |\mathbf{e}_L \cdot \mathbf{e}_R|^2 = \left| \left( 1/\sqrt{2}, -i/\sqrt{2}, 0 \right) \cdot \left( 1/\sqrt{2}, -i/\sqrt{2}, 0 \right) \right|^2 = \left| \frac{1}{2} - \frac{1}{2} \right|^2 = 0
\end{align*}
\]

**Irrelevant point:** Sometimes, you can find the direction of \( \mathbf{k} \) from the polarization vector, sometimes not. If \( \mathbf{e}_R \) and \( \mathbf{e}_I \) point in different directions, then because they are both perpendicular to \( \mathbf{k} \), \( \mathbf{e}_R \times \mathbf{e}_I \) points along \( \mathbf{k} \). But if \( \mathbf{e}_R \) and \( \mathbf{e}_I \) are parallel, or if \( \mathbf{e} = 0 \), then \( \mathbf{e}_R \times \mathbf{e}_I = 0 \), and you can’t tell anything. So far as I know, this fact is of no use at all.

**TBS.** Angular momentum of elliptically polarized waves.

**Extension to other waves:** For things like gravity waves, the field which varies sinusoidally is a rank-2 tensor, which can be written as a 2D matrix (below, left):

\[
E_{\mu\nu}(t, x) = \text{Re} \left[ \begin{array}{cccc}
E_{tt} & E_{tx} & E_{ty} & E_{tz} \\
E_{xt} & E_{xx} & E_{yx} & E_{xz} \\
E_{yt} & E_{yx} & E_{yy} & E_{yz} \\
E_{zt} & E_{zx} & E_{zy} & E_{zz}
\end{array} \right] e^{-i\omega t}
\]

\[
E_{z}(t, x) = \frac{1}{\sqrt{2}} \text{Re} \left[ \begin{array}{cccc}
0 & 0 & 0 & 0 \\
0 & 1 & 0 & 0 \\
0 & 0 & 1 & 0 \\
0 & 0 & 0 & 0
\end{array} \right] e^{-i\omega t}
\]

Above right is \( \tilde{\mathbf{e}}_z \), the polarization tensor for “+” polarization propagating in the \( z \)-direction.

**Sleepy Hollow: The Legend of the Headless Vector**

Linear polarization direction is often called a “headless vector,” because often only the plane of polarization is important (or known). For example, if we rotate a linear polarizer by 180°, it’s behavior is not changed. Similarly, if we rotate the light incident on a polarizer by 180°, there is no change in the output light. However, these experiments do not explore the full “headedness” of polarization. We now show that polarization is, in fact, a “headful” vector, whose sense can be measured with interference. This
is reasonable, because the polarization is fully described by the direction of the \( E \)-field, at a given reference time, and the \( E \)-field is a headful vector.

Recall every vector lies along an axis, and that the sense of a vector is simply which way the vector points along the axis. In some cases, including most cases of polarization, we know the axis of the \( E \)-field, but not its sense. Also, recall that the polarization vector specifies the complete (“headful”) direction and phase of the \( E \)-field for a propagating wave.

Interference occurs between headful \( E \)-field vectors, including their sense. To illustrate, suppose we superpose two equal-amplitude linearly polarized EM waves, with polarization vectors \( \vec{e}_1 \) and \( \vec{e}_2 \). Then at the point of superposition, we get:

\[
\begin{align*}
\vec{E}_1(t) &= \text{Re} \left\{ \vec{E}_0 \vec{e}_1 e^{-i\omega t} \right\}, \\
\vec{E}_2(t) &= \text{Re} \left\{ \vec{E}_0 \vec{e}_2 e^{-i\omega t} \right\}, \\
\vec{E}_1 + \vec{E}_2 &= \text{Re} \left\{ \vec{E}_0 \left( \vec{e}_1 + \vec{e}_2 \right) e^{-i\omega t} \right\}.
\end{align*}
\]

Suppose that \( \vec{e}_1 \) and \( \vec{e}_2 \) are real and parallel: \( \vec{e}_1 = \vec{e}_2 \). Then we get constructive interference. Now suppose we rotate \( \vec{e}_2 \) in space the \( 2^{nd} \) wave about its propagation direction by \( 180^\circ \). This is not a phase-shift; there is no time delay involved. After rotation, the polarization vector gets negated: \( \vec{e}_2 \to -\vec{e}_1 \). The interference is now destructive:

\[
\vec{E}_1 + \vec{E}_2 = \text{Re} \left\{ \vec{E}_0 \left( \vec{e}_1 + \vec{e}_2 \right) e^{-i\omega t} \right\} = 0.
\]

Therefore, we see that \( \vec{e}_1 \) and \( -\vec{e}_1 \) (or \( \vec{e}_2 \) and \( -\vec{e}_2 \)) are different polarizations. One is the negative of the other. If \( \vec{e}_2 = (\vec{e}_0, 0, 0) \) then it is horizontal polarization. Then \( -\vec{e}_2 \) is also horizontal polarization, but it’s a different “horizontal:” it is the negative of \( \vec{e}_1 \)’s “horizontal.”

Thus the polarization vector describes not only the plane of polarization, but the sense of the \( E \)-field in that plane. For linear polarization, you can write the polarization as a unit vector parallel to the \( E \)-field. This is a simple case of a polarization vector.

Of course, the sense reverses every half-period, so the sense of the polarization vector is defined at some reference time. In fact, the polarization vector describes the full \( 360^\circ \) phase (in time) of the \( E \)-field, and phase is always relative to some reference time (or phase), taken as zero. The phenomenon of constructive or destructive interference is, as always, independent of our reference time (or phase).

**Summary:** Some experiments are not sensitive to the sense of the polarization, and therefore measure only the plane of polarization. They often treat such a polarization measurement as a headless vector, which is good enough for some applications. However, we have shown that interference reveals that polarization is a headful vector, even if the sense is unknown.

---

**Poynting Vector For Linear Polarization**

The Poynting vector describes the power density (= energy flux) of the propagating radiation: watts/m\(^2\). It points in the direction of power flow, and is proportional to \( E \) and \( B \):

\[
\vec{S}(t, \mathbf{r}) = \vec{E}(t, \mathbf{r}) \times \vec{H}(t, \mathbf{r}) \quad \text{SI units: } (\text{V/m})(\text{A/m}) = \text{W/m}^2.
\]

Since \( H \propto E \), and perpendicular to it, if either \( E \) or \( H \) is sinusoidal, then both are sinusoidal:

\[
\vec{S}(t) = EH \sin^2 \omega t \quad \text{where } \quad E, H = \text{amplitudes}.
\]

Quite often, we are more interested in the time-averaged power density, rather than the instantaneous power density. Then we use the fact that \( \langle \sin^2 \rangle = \frac{1}{2} \), over the long term (or an integral number of quarter-cycles):

\[
\left\langle \vec{S} \right\rangle = \frac{1}{2} EH \left( \hat{\vec{E}} \times \hat{\vec{H}} \right), \quad \text{where } \quad \hat{\vec{E}} = \frac{\vec{E}}{E}, \quad \hat{\vec{H}} = \frac{\vec{H}}{H} \quad \text{are unit vectors in direction of } \vec{E} \text{ and } \vec{H}.
\]
Beware of Solenoidal Poynting Vectors

In the radiation formula for a point particle [Jac ?? & Gri ??], there is a term called the “velocity term” or “velocity field,” because it is proportional to the particle’s velocity:

\[ \text{radiation formula here ??} \]

This term seems to contribute to a Poynting vector, yet is said not to contribute to radiation. Some references’ explanations are misleading: they say that the electromagnetic intensity due to the velocity term falls off as $1/r^4$, and so becomes insignificant compared to the $1/r^2$ “radiation” term. While the velocity term does drop as $1/r^4$, this is not the crux of the matter. The Poynting vector describes the power density (= energy flux density) of the propagating radiation: watts/m$^2$. Regardless of how the magnitude of the Poynting vector falls off, a radial component to it represents real energy flow. In the far field, we can imagine thick spherical shells around the radiation source. Since the shells are in vacuum (or non-absorbing material), the power (energy/time) flowing into any shell (from the source inside) must equal the power flowing out of that shell. Since the shell surface area increases as $r^2$, the power density must decrease as $1/r^2$, so that the product of the two, the total power, remains constant. This is just conservation of energy. Therefore:

**Conservation of energy requires that all radially directed power flows fall off exactly as $1/r^2$.**

There cannot be any power flow that falls off as $1/r^4$, because it would violate conservation of energy. So what of the velocity term? The real explanation is that the Poynting vector resulting from the velocity term is solenoidal, i.e. the lines of power flow close on themselves. No power flows out; it’s almost as if the power just flows around in circles. The radial component of the velocity term Poynting vector is 0:

\[ \mathbf{S}_{\text{velocity}} \cdot \hat{r} = 0 \]  

(velocity term Poynting vector).

Velocity fields have no outward power flow. The rate of fall off ($1/r^4$ or anything else) is irrelevant. We can see this by considering the well-known E-field of a particle at constant velocity. This field emanates from a single point in space called the “present position.” That means the E-field everywhere points radially outward. It is also well-known that only the tangential component of E contributes to radiation, which you can readily see because only the perpendicular component of E contributes to the radial component of the Poynting vector (diagram??). The velocity field has no tangential component of E, and therefore no radiation. The Poynting vector field-lines make circles around the present position, and are completely solenoidal. We can compute the radiated power out of an arbitrary surface (not necessarily a sphere, nor centered anywhere in particular) from:

\[ P_{\text{rad}} = \oiint S_n \, da. \]

While there may be patches where the velocity terms have an outward component, their solenoidal nature insures that any such patch is exactly canceled by an inward component somewhere else.

In short, velocity terms don’t radiate because their Poynting vectors go in closed loops.

TBS: Poynting Vector For Arbitrary Polarization

Wave Packets

Often, one sends finite-time “messages” or signals with electromagnetic waves (or some other linear medium: wires, water, air, etc.). Such a signal can be decomposed into a sum of sinusoidal frequency components (i.e., the Fourier Transform), each frequency component with its own amplitude and phase. Since many propagation phenomena are frequency dependent, it is important to ask: What range of frequencies are needed to add together to form a finite time signal? We will argue that a continuous “band” of frequencies, of finite upper and lower bound, is sufficient to construct a finite time signal. The difference between the upper and lower frequency in a signal is called the **bandwidth** (BW). (This term is much abused; in particular, it is frequently used (incorrectly) to mean “data capacity.”) We give a hand-waving argument for a crude estimate of the bandwidth of a finite-time signal. When viewed as a sum of sinusoidal waves, such a signal is called a **wave-packet**.
Figure 3.8 (Left) A wave-packet as a function of time. (Right) Frequency components of a wave-packet.

Note that when you transmit time-signals into a propagating medium, their extension in time becomes also an extension in space (e.g., wavelength). The following argument applies equally well to a signal at a point in space that varies over time, and to a signal frozen in time that varies in space. Any reference to frequency can be equivalently thought of as time-frequency ($\omega$), or spatial-frequency (wave-number, $k$).

When adding up lots of sinusoids, we consider the instantaneous phase of each component at some point (in either time or space). If there is a wide range of phases, spread over most of $[0, 2\pi)$, then the sinusoids will have lots of positive and negative components, and will add up to approximately zero. To have a significant contribution to our signal, most of the instantaneous phases need to be within a radian or two of each other. And this has to be true throughout the width (in time or space) of the wave packet.

Since the signal comprises different frequencies, the instantaneous phase difference between frequency components varies over time (or space). Components “in-phase” at one point will be “out-of-phase” at a distant point. (Below) the low- and high-frequency components (red and blue) start out in-phase, but drift to opposite phase over some time.

Figure 3.9 Wave components of different frequencies drift out of phase over time (or space).

The rate at which the phases drift apart is exactly the frequency difference between the two components:

$$\theta_1 = \omega_1 t + \phi_1, \quad \theta_2 = \omega_2 t + \phi_2$$

$$\Delta \theta = \theta_1 - \theta_2 = \omega_1 t + \phi_1 - \omega_2 t - \phi_2 = (\omega_1 - \omega_2) t + \text{const} = (\Delta \omega) t + \text{const}, \quad \text{or} \quad \Delta \theta = (\Delta k) z + \text{const}$$

The constant term is not important, and simply depends on the shape of the signal and the origin of our $t$ (or $z$) coordinate. What matters is that as $t$ (or $z$) increases, so does the phase difference. Choose the coordinate origin to be the start of the wave-packet; then the width of the packet is $\Delta t$ (or $\Delta z$). Clearly, the largest $\Delta \theta$ occurs when the frequency difference is maximum, i.e., between $\omega_{lower}$ and $\omega_{upper}$:

$$\Delta \theta_{\text{max}} = (\omega_{upper} - \omega_{lower}) \Delta t = (BW) \Delta t, \quad \text{or} \quad \Delta \theta_{\text{max}} = (k_{upper} - k_{lower}) \Delta z = (BW) \Delta z.$$

Since $\Delta \theta$ must remain within 1 or 2 radians over the width of the signal, we have

$$\Delta \theta \leq 1 \text{ or } 2 \text{ rad} = (BW) \Delta t \Rightarrow \quad BW \leq \frac{1 \text{ or } 2 \text{ rad}}{\Delta t}, \quad \text{or} \quad BW \leq \frac{1 \text{ or } 2 \text{ rad}}{\Delta z}.$$
A more detailed analysis yields a theorem which states: a signal of finite time must (strictly speaking) have infinite bandwidth, and conversely, a signal of finite bandwidth must extend over infinite time. However, one can have an approximately finite-time and finite-bandwidth signal (for all practical purposes), because the amplitudes outside a small interval are negligibly small.

**Phase Velocity and Group Velocity**

A hugely important concept now arises: what is the propagation speed of the wave packet? You might be tempted to say, “the same as the propagation speed of a single frequency.” That’s true if all frequencies propagate at the same speed. The speed of a single sinusoid is called the **phase velocity**, and can be seen from Figure 3.10.

In the figure, we have $4\pi$ rad/s, and $\pi$ rad/m. The wave must then travel with:

$$v_{\text{phase}} = \frac{\omega}{k} = \frac{4\pi \text{ rad/s}}{\pi \text{ rad/m}} = 4 \text{ m/s}.$$ 

If all frequencies propagate at the same speed, then $\omega = v_{\text{phase}}k$ (Figure 3.10, right).

However, in most linear media, the propagation speed of a sinusoid **depends** on its frequency. Since a wave-packet is made up of many frequencies, which frequency should we use to determine its speed? You might guess the center frequency. Turns out, though, that what really matters is just how the speed **varies** with frequency. The wave-packet, or “group” of waves travels with a speed (**group-velocity**):

$$v_{\text{group}} = \frac{d\omega}{dk}. \quad \text{Recall the phase velocity } v_{\text{phase}} = \frac{\omega}{k}.$$ 

Actually, since each component travels at a different speed, the wave packet does not exactly travel “as a group.” It gets “smeared” over time, because the relative phases of its components get misaligned. This smearing is called **dispersion**, and in some cases, is a serious limitation on a communication system. More on phase and group velocity in “Waveguides,” later.

**Vector Potentials I Have Known**

Many E&M questions are well-suited to using the magnetic vector potential, $A(\mathbf{r})$, for the magnetic fields. Given a current distribution, we can (in principle) find the vector potential from:

$$A(\mathbf{r}) = \frac{\mu_0}{4\pi} \int \frac{J(\mathbf{r}')}{|\mathbf{r} - \mathbf{r}'|} \, d^3r'. \quad \text{[Jac 5.32 p181]}.$$ 

However, many problems start with a given magnetic field (rather than current distribution), but our analysis tools use the vector potential. How can we find the vector potential for a given magnetic field? There is no explicit “inverse curl,” but usually it is fairly easy to do by inspecting the formula for curl in the
coordinates of interest. For example, for a constant \( B \)-field, \( \mathbf{B} = \nabla \times \mathbf{A} = B\hat{z} \) and rectangular coordinates (chosen from other symmetries of the problem), we look at the formula for curl in those coordinates:

\[
\nabla \times \mathbf{A} = \left( \frac{\partial A_x}{\partial y} - \frac{\partial A_y}{\partial z} \right) \hat{x} + \left( \frac{\partial A_y}{\partial z} - \frac{\partial A_z}{\partial x} \right) \hat{y} + \left( \frac{\partial A_z}{\partial x} - \frac{\partial A_x}{\partial y} \right) \hat{z}.
\]

We are given only a \( B_z \) component, so we look at the coefficient of \( \hat{z} \) above. We see that we can get it from either a \( \partial A_y/\partial x \), or a \( \partial A_x/\partial y \) term. Both terms are simple, so we choose one, and set:

\[
\frac{\partial A_y}{\partial x} = B_z \quad \Rightarrow \quad A_y(x) = \int_0^r dx B_z = B_z x.
\]

If other aspects of the problem favored an \( A_x \) instead, we could choose the other term in the curl. Note that the curl operator is linear, so if there are more components of \( B \), we can build them up from \( A \) terms, and sum them for the final \( A \).

Suppose we have a constant \( B_z \) in cylindrical coordinates,

\[
\nabla \times \mathbf{A} = \left( \frac{1}{r} \frac{\partial A_z}{\partial \phi} - \frac{\partial A_\phi}{\partial z} \right) \hat{r} + \left( \frac{\partial A_y}{\partial z} - \frac{\partial A_z}{\partial r} \right) \hat{\phi} + \left( \frac{r}{\partial r} \frac{\partial A_\phi}{\partial \phi} - \frac{\partial A_r}{\partial \phi} \right) \hat{\phi}
\]

Again we see two terms to choose from: \( \frac{1}{r} \frac{\partial}{\partial \phi} (rA_\phi) \) and \( \frac{1}{r} \frac{\partial}{\partial \phi} (rA_\phi) \). If the problem has axial symmetry, then the 2\(^{nd} \) term must be zero. Also, the 2\(^{nd} \) term cannot have a constant derivative, and be single valued at \( \phi = 0 = 2\pi \). So we choose the first term, and set

\[
\frac{1}{r} \frac{\partial}{\partial \phi} (rA_\phi) = B_z
\]

\[
\frac{1}{r} \frac{\partial}{\partial \phi} (rA_\phi) = B_z \quad \Rightarrow \quad A_\phi = B_z/r.
\]

Spherical coordinates don’t come up all that often in magnetics problems, and are a little more complicated, but we can use the same method.

\[
\nabla \times \mathbf{A} = \frac{1}{r \sin \theta} \left( \frac{\partial}{\partial \theta} (\sin \theta A_\phi) - \frac{\partial A_\theta}{\partial \phi} \right) \hat{r} + \left( \frac{1}{r \sin \theta} \frac{\partial A_y}{\partial \phi} - \frac{1}{r} \frac{\partial}{\partial r} (rA_\phi) \right) \hat{\phi} + \left( \frac{\partial}{\partial r} (rA_\phi) - \frac{\partial A_\phi}{\partial \theta} \right) \hat{\phi}
\]

For a constant \( B_z \), we must first decompose \( \mathbf{z} \) into its \( r \) and \( \theta \) components:

\[
\hat{z} = \cos \theta \hat{r} - \sin \theta \hat{\theta} \Rightarrow \frac{1}{r \sin \theta} \frac{\partial}{\partial \theta} (\sin \theta A_\phi) - \frac{\partial A_\theta}{\partial \phi} = \cos \theta
\]

\[
\frac{1}{r \sin \theta} \frac{\partial A_y}{\partial \phi} - \frac{1}{r} \frac{\partial}{\partial r} (rA_\phi) = -\sin \theta
\]

Choose: \( A_\theta = 0 \) \( \Rightarrow \frac{1}{r \sin \theta} \frac{\partial}{\partial \theta} (\sin \theta A_\phi) = \cos \theta
\]

\[
\sin \theta A_\phi = r \int d\theta \cos \theta \sin \theta = \frac{r}{2} \sin^2 \theta \quad \Rightarrow \quad A_\phi = \frac{r}{2} \sin \theta
\]

Plug into the 2\(^{nd} \) equation from \( \hat{z} \):
Solving Laplace’s Equation (Boundary Value Problems)

Although these are generally covered in standard texts, we take a different approach from any we have seen. Note that boundary value problems (BVPs) are essential before moving on to waveguides. We describe solutions to Laplace’s equation in eigenfunction-like terms. This has a huge advantage when moving on to waveguides, which are just big eigenfunction boundary value problems. It also is familiar to most students, because eigenfunctions appear everywhere. It can even serve as a simple introduction to eigenfunctions.

We consider both 2-D and 3-D BVPs, in all common coordinate systems: rectangular, polar, cylindrical, and spherical. All problems start from Laplace’s equation:

\[ \nabla^2 \Phi = 0, \]

plus some boundary conditions on \( \Phi \). In all common coordinate systems, the nontrivial solutions arise from separations of variables. Separation of variables hopes that we can write the solution as a product of functions, each of a single coordinate, such as:

\[ \Phi(x, y, z) = X(x)Y(y)Z(z), \quad \Phi(r, \phi, z) = R(r)Q(\phi)Z(z), \quad \text{or} \quad \Phi(r, \phi, \theta) = R(r)Q(\phi)P(\theta). \]

We restrict ourselves to real-valued functions and parameters, since they are of most interest to E&M. [However, quantum mechanics uses complex functions which are simple linear combinations of the real functions we use here.] Finally, note that Laplace’s equation is linear, so that the weighted sum (linear superposition) of any number of solutions is also a solution. We use this property extensively in what follows.

We cover the various cases in order of difficulty. We do not consider orthogonal function expansions for matching boundary conditions; see regular texts [refs??].

Two-D Laplace Solutions

Recall that the 2-D Laplacian operator is:

\[ \nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2}, \quad \text{or} \quad \nabla^2 = \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2}. \]

Two-D Laplace Solutions in Rectangular Coordinates

We start with the simplest solution to \( \nabla^2 \Phi = 0 \), which is completely omitted from Jackson [Jac ch 2]:

\[ \Phi(x, y) = Ax + By + C. \]

By inspection, this clearly satisfies Laplace’s equation. Both terms of the Laplacian operator are zero. It turns out, though, that this solution doesn’t come up very often in E&M problems, because it doesn’t satisfy most physical boundary conditions. Nonetheless, it is a perfectly valid solution, and should be considered in some cases. [It comes up in some 1D quantum mechanical scattering problems.]

More interestingly, we consider separation of variables, which requires:

\[ \Phi(x, y) = X(x)Y(y) . \]

Consider the two terms of the Laplacian in rectangular coordinates, i.e., the partial derivatives w.r.t. \( x \) and \( y \):

\[ \frac{1}{r \sin \theta} \frac{\partial A_r}{\partial \phi} - \frac{1}{\sin \theta} \frac{\partial}{\partial r} \left( \frac{r}{2} \sin \theta \right) \]

\[ = -\sin \theta \]

\[ \frac{1}{r \sin \theta} \frac{\partial A_r}{\partial \phi} - \sin \theta = -\sin \theta, \quad \frac{\partial A_r}{\partial \phi} = 0 \Rightarrow A_r = \text{const}, \ \text{choose} \ A_r = 0 \]
\[ \frac{\partial^2 \Phi}{\partial^2 x} = \left( \frac{\partial^2 X}{\partial^2 x} \right) Y(y), \quad \frac{\partial^2 \Phi}{\partial^2 y} = X(x) \frac{\partial^2 Y}{\partial^2 y} \Rightarrow \left( \frac{\partial^2 X}{\partial^2 x} \right) Y(y) + X(x) \frac{\partial^2 Y}{\partial^2 y} = 0. \]

If we could find \( X(x) \) and \( Y(y) \) with these properties, we’d be set:

\[ \frac{\partial^2 X}{\partial x^2} = kX(x), \quad \text{and} \quad \frac{\partial^2 Y}{\partial y^2} = -kY(y), \quad \text{then} \]

\[ \nabla^2 \Phi = kX(x)Y(y) - kX(x)Y(y) = (k - k)X(x)Y(y) = 0. \]

This implies that \( X(x) \) and \( Y(y) \) are eigenfunctions of the 2nd derivative operator. In rectangular coordinates, we can easily find such a solution. Recall that \( \sinh(x) \) and \( \cosh(x) \) (and equivalently \( \exp(x) \) and \( \exp(-x) \)) satisfy this, and their eigenvalues are always positive. Also, \( \sin(x) \) and \( \cos(x) \) satisfy this, and their eigenvalues are always negative. The solutions are therefore:

\[
\begin{align*}
X(x) &= Ae^{\alpha x} + Be^{-\alpha x}, \\
Y(y) &= C \sin \alpha y + D \cos \alpha y,
\end{align*}
\]

\[ \frac{\partial^2 X}{\partial x^2} = \alpha^2 \left( Ae^{\alpha x} + Be^{-\alpha x} \right) = \alpha^2 X(x), \quad \text{and} \]

\[ \frac{\partial^2 Y}{\partial y^2} = -\alpha^2 \left( C \sin \alpha y + D \cos \alpha y \right) = -\alpha^2 Y(y) \]

OR

\[
\begin{align*}
X(x) &= C \sin \alpha x + D \cos \alpha x, \\
Y(y) &= Ae^{\alpha y} + Be^{-\alpha y},
\end{align*}
\]

\[ \frac{\partial^2 X}{\partial x^2} = -\alpha^2 \left( C \sin \alpha x + D \cos \alpha x \right) = -\alpha^2 X(x), \quad \text{and} \]

\[ \frac{\partial^2 Y}{\partial y^2} = \alpha^2 \left( Ae^{\alpha y} + Be^{-\alpha y} \right) = \alpha^2 Y(y) \]

\[ \Phi(x, y) = X(x)Y(y) \]

As always, the boundary conditions determine the coefficients \( A, B, C, \) and \( D. \)

Note that

\[ \sinh(x) = \frac{1}{2} e^x - \frac{1}{2} e^{-x}, \quad \cosh(x) = \frac{1}{2} e^x + \frac{1}{2} e^{-x}, \quad \text{and} \]

\[ A \sin(x) + B \cos(x) = E \cos(x + \beta), \quad \text{where} \quad E = \sqrt{A^2 + B^2}, \quad \beta = \tan^{-1} \left( \frac{B}{A} \right). \]

\( \sinh(x) \) and \( \cosh(x) \) and \( \exp(x) \) and \( \exp(-x) \) are linear combinations of each other, and therefore equally good solutions. However, \( \sinh(x) \) and \( \cosh(x) \) are more often convenient in E&M problems because \( \sinh \) has a zero, and \( \cosh \) has a zero derivative, which are useful for matching common boundary conditions. \( \exp(\pm x) \) have no zeros, and no zero derivatives. Also, \( E \cos(x + \beta) \) is equivalent to \( A \sin(x) + B \cos(x) \). So we can conveniently write:
\[
\begin{align*}
X(x) &= A \sinh(\alpha x) + B \cosh(\alpha x), \\
\frac{\partial^2 X}{\partial x^2} &= \alpha^2 (A \sinh(\alpha x) + B \cosh(\alpha x)) = \alpha^2 X(x), \quad \text{and} \quad \\
Y(y) &= E \cos(\alpha y + \beta), \\
\frac{\partial^2 Y}{\partial y^2} &= -\alpha^2 E \cos(\alpha y + \beta) = -\alpha^2 Y(y)
\end{align*}
\]

OR
\[
\begin{align*}
X(x) &= E \cos(\alpha x + \beta), \\
\frac{\partial^2 X}{\partial x^2} &= -\alpha^2 E \cos(\alpha x + \beta) = -\alpha^2 X(x), \quad \text{and} \quad \\
Y(y) &= A \sinh(\alpha y) + B \cosh(\alpha y), \\
\frac{\partial^2 Y}{\partial y^2} &= \alpha^2 (A \sinh(\alpha y) + B \cosh(\alpha y)) = \alpha^2 Y(y)
\end{align*}
\]
\[
\Phi(x, y) = X(x)Y(y).
\]

As always, the boundary conditions determine \( A, B, E, \) and \( \beta \) in this alternate form.

More on boundary conditions later.

**Two-D Laplace Solutions in Polar Coordinates**

In polar coordinates, \((r, \phi)\), we might try the same trick for separation of variables: eigenfunctions of the two Laplacian operator terms, with eigenvalues that sum to zero:

\[
\nabla^2 = \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} \quad \text{Let} \quad \Phi(r, \phi) = R(r)Q(\phi). \quad \text{Solve}
\]

\[
\frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) R(r) = kR(r), \quad \quad \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = -kQ(\phi)
\]

We see immediately that this won’t work, because of the \(1/r^2\) factor in front of the \(Q(\phi)\) term, and \(Q(\phi)\) has no \(r\) in it, by definition. But we don’t really need *eigenfunctions* of the Laplacian terms; we just need a function \(f(r, \phi)\) such that:

\[
\frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) R(r) = f(r)R(r), \quad \quad \text{and} \quad \\
\frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = -f(r)Q(\phi). \quad \quad \text{Then:}
\]

\[
\nabla^2 \Phi = f(r)R(r)Q(\phi) - R(r)f(r)Q(\phi) = 0
\]

Given that the 2nd term of the Laplacian has \(1/r^2\) before it, the simplest choice for \(f(r)\) that we could hope for is:
Let \( f(r) = \frac{k}{r^2} \).

Then \( R(r) = Ar^v + Br^{-v} \) \quad \Rightarrow \quad \frac{1}{r} \frac{\partial}{\partial r} \left( \frac{r}{\partial r} \right) R(r) = \frac{1}{r} \frac{\partial}{\partial r} \left( r \left( vAr^{-v-1} - vBr^{-v-1} \right) \right)
\[
= \frac{1}{r} \frac{\partial}{\partial r} \left( vAr^v - vBr^{-v} \right) = \frac{1}{r} \left( v^2 Ar^{-v} + v^2 Br^{-v-1} \right) = \frac{v^2}{r^2} R(r)
\]

and \( Q(\phi) = C \sin v\phi + D \cos v\phi \) \quad \Rightarrow \quad \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = \frac{1}{r^2} \left( -v^2 C \sin v\phi - v^2 D \cos v\phi \right) = -\frac{v^2}{r^2} Q(\phi)
\]

So \( f(r) = \frac{v^2}{r^2} \),

and we have found our solution to Laplace’s equation in 2-D polar coordinates. [It happens now that \( Q(\phi) \) is an eigenfunction of \( \frac{\partial^2}{\partial \phi^2} \), but \( R(r) \) is not an eigenfunction of anything.] Note that, depending on the boundary conditions, \( v \) may or may not be an integer. [Jac sec 2.11 p75] covers this well. In particular, if \( \phi \) goes completely around the origin, then \( v \) must be an integer, because it must be that \( Q(\nu(2\pi + \phi)) = Q(\nu\phi) \), so that \( Q \) is single-valued at every point. The solutions are thus

\[
R(r) = Ar^v + Br^{-v}, \quad Q(\phi) = C \sin v\phi + D \cos v\phi, \quad \Phi(r, \phi) = R(r)Q(\phi).
\]

Note that if the domain of \( \Phi \) includes the origin, then \( R(r) = Ar^v \), because \( Br^{-v} \) blows up there.

There is one more case: \( v = 0 \). If this is allowed by boundary conditions and the domain of \( \Phi \), then the two Laplacian terms are both zero. There are separate solutions for this case:

\[
\frac{1}{r} \frac{\partial}{\partial r} \left( \frac{r}{\partial r} \right) R(r) = 0, \quad \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = 0
\]

\[
R(r) = B \ln \frac{r}{A} \quad \text{(sometimes written} \quad R(r) = K + B \ln r) \quad \Rightarrow \quad \frac{1}{r} \frac{\partial}{\partial r} \left( \frac{r}{\partial r} \right) R(r) = \frac{1}{r} \frac{\partial}{\partial r} \left( \frac{B}{r} \right) = \frac{1}{r} \frac{\partial}{\partial r} B = 0
\]

and \( Q(\phi) = C + D\phi \) \quad \Rightarrow \quad \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = 0
\]

\[
\Phi(r, \phi) = R(r)Q(\phi) = \left( B \ln \frac{r}{A} \right) \left( C + D\phi \right)
\]

Note that this possibility exists only if the domain of \( \Phi \) excludes the origin, because \( \ln r \) blows up there. Also note that if the domain of \( \Phi \) allows \( \phi \) to go all the way around the origin, then \( D = 0 \), so that \( Q(\phi) \) is single valued, and hence \( Q(\phi) = 1 \) (constant).

In this case of the domain surrounding but excluding the origin, we get the most general solution by combining the \((v = 0, D = 0)\) case with the other integer \( v \) [Jac 2.71 p77]:

\[
\Phi(r, \phi) = B_0 \ln \frac{r}{A_0} + \sum_{\nu=1}^{\infty} \left( A_\nu r^\nu + B_\nu r^{-\nu} \right) \left( C_\nu \sin \nu\phi + D_\nu \cos \nu\phi \right).
\]

7/7/2017 16:32  Copyright 2002 - 2017 Eric L. Michelsen. All rights reserved.  Page 29 of 75
Three-D Laplace Solutions

**Three-D Laplace Solutions in Rectangular Coordinates**

First, we extend the almost-trivial solution, which satisfies Laplace’s equation by inspection:

\[ \nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2}, \quad \Phi(x, y, z) = Ax + By + Cz + D. \]

All 3 Laplacian terms are zero. Again, it doesn’t come up much, but some references don’t even mention it.

For less trivial solutions, we extend the separation of variables approach from the 2D rectangular coordinate case:

\[ \Phi(x, y, z) = X(x)Y(y)Z(z) \]

\[ \frac{\partial^2 X}{\partial x^2} = aX(x), \quad \frac{\partial^2 Y}{\partial y^2} = bY(y), \quad \frac{\partial^2 Z}{\partial z^2} = cZ(z), \quad \text{where} \quad a + b + c = 0. \]

If any coordinate, say \( z \), can be separated into the form \( Cz + D \), its eigenvalue is zero, and we revert to the 2D solutions for the remaining two coordinates. In the 2D case, we had only two terms, so the eigenvalues had to be negatives of each other, so that they would sum to zero, and satisfy Laplace’s equation. Recall that the eigenvalues of \( \sinh(x) \) and \( \cosh(x) \) are always positive, and the eigenvalues of \( \sin(x) \) and \( \cos(x) \) are always negative. This meant that one coordinate had to be \( A \sinh(\cdot) + B \cosh(\cdot) \) (or equivalently \( \exp(\pm \cdot) \)), and the other had to be \( C \sin(\cdot) + D \cos(\cdot) \).

In the 3-D case, there are more possibilities, but all we require is that the 3 eigenvalues sum to zero:

\[ a + b + c = 0. \]

This means at least one function must be \( \sinh(\cdot) \) or \( \cosh(\cdot) \) (or equivalently \( \exp(\pm \cdot) \)), at least one function must be \( \sin \) or \( \cos \), and the last function can be either, subject to the constraint that the eigenvalues of all 3 functions sum to zero. Some example solutions are:

\[ X(x) = \sin 2x, \quad Y(y) = \cos 3y, \quad Z(z) = \sinh \left( \sqrt{1/3} z \right) \quad \left( -2^2 - 3^2 + 13 = 0 \right) \]
\[ X(x) = \cosh(3x), \quad Y(y) = \sinh(4y), \quad Z(z) = \cos(5z) \quad \left( 3^2 + 4^2 - 5^2 = 0 \right) \]
\[ X(x) = e^{5x}, \quad Y(y) = e^{-12y}, \quad Z(z) = 2 \sin(13z) + 7 \cos(13z) \quad \left( 5^2 + 12^2 - 13^2 = 0 \right) \]

**Three-D Laplace Solutions in Cylindrical Coordinates**

In cylindrical coordinates, if the \( z \) coordinate can be separated into the form \( Cz + D \), its eigenvalue is zero, and we revert to the 2D solutions for polar coordinates. This is rare.

Barring that, we now must introduce a slight variation on our separation of variables theme: the Bessel functions. We start with the standard approach of separation of variables:

\[ \nabla^2 = \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} + \frac{\partial^2}{\partial z^2}, \quad \Phi(r, \phi, z) = R(r)Q(\phi)Z(z). \quad (3.1) \]

The sum of the above 3 Laplacian terms must be zero. From our 2D work, we have a \( Q(\phi) = \sin \) or \( \cos(\nu \phi) \) solution that can offset an \( R(r)/r^2 \) term, and a \( Z(z) \) solution that can offset either a positive or negative constant term \( [Z(z) = \sinh \text{ or } \cosh(kz) \text{ or } Z(z) = \sin \text{ or } \cos(kz)] \):
\[
\frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = -\frac{\nu^2}{r^2} Q(\phi),
\]
(3.2)

\[
\frac{\partial^2 Z}{\partial z^2} Z(z) = k^2 Z(z).
\]

Wouldn’t it be nice if we had an \( R(r) \) that would produce the negative of the sum of those \( Q(\phi) \) and \( Z(z) \) terms? That is:

\[
\frac{1}{r} \frac{\partial}{\partial r} \left( \frac{\partial}{\partial r} \right) R(r) = \left( -k^2 + \frac{\nu^2}{r^2} \right) R(r)
\]
(3.3)

This equation is singular at \( r = 0 \), which is often a physically valid radius. The singularity is easily removed by defining \( R(r) \) as the limit:

\[
R(0) = \lim_{r \to 0} R(r).
\]

**Case 1:** We call the eigenvalue of \( Z(z) \) “\( k^2 \)”, and it is positive. \( Z(z) = \sinh \text{ or } \cosh(kz) \).

We don’t have any pre-existing functions that satisfy the \( R(r) \) equation, so we do the obvious: we make one! Using power series solutions and some fancy math (that we don’t need to know here), mathematicians define a family of functions which satisfy the above \( R(r) \) equation for \( k^2 > 0 \), and call them Bessel functions. Every value of \( \nu \) gets its own Bessel function, named \( J_\nu(kr) \). Recall that \( \nu \) can be any real number, so there is an uncountably infinite set of Bessel functions. Bessel functions have lots of magic properties (that we don’t need to know here). We don’t need a separate Bessel function for each value of \( k \), because a simple change of variable eliminates \( k \) (\( w = kr \)) [I need to clarify this??]:

Let \( w = kr \), \( r = \frac{w}{k} \). Then \( \frac{\partial}{\partial r} \to k \frac{\partial}{\partial w} \), and

\[
\frac{1}{r} \frac{\partial}{\partial r} \left( \frac{\partial}{\partial r} \right) R(r) = \left( -k^2 + \frac{\nu^2}{r^2} \right) R(r) \to \frac{w}{k} \frac{k}{w} \frac{\partial}{\partial w} \frac{w}{k} \frac{\partial}{\partial w} \frac{w}{k} \frac{\partial}{\partial w} R \left( \frac{w}{k} \right) = \left( -k^2 + \frac{k^2 \nu^2}{w^2} \right) R \left( \frac{w}{k} \right)
\]

Define \( J_\nu(w) = R \left( \frac{w}{k} \right) \). Then

\[
\frac{1}{w} \frac{\partial}{\partial w} \left( w \frac{\partial}{\partial w} \right) J_\nu(w) = \left( -1 + \frac{\nu^2}{w^2} \right) J_\nu(w)
\]

Then \( R \left( \frac{w}{k} \right) = J_\nu(w) \) \( \Rightarrow \) \( R(r) = J_\nu(kr) \) \( \Rightarrow \)

\[
\Phi(r, \phi, z) = J_\nu(\nu) \left( C \sin \nu \phi + D \cos \nu \phi \right) \left( E \sinh k z + F \cosh k z \right)
\]

Notice that since the eigenvalue of \( Z(z) > 0 \) (\( k^2 > 0 \)), \( Z(z) = \sinh(kz) \) or \( \cosh(kz) \). (We’ll get to the \( Z(z) = \sin \) or \( \cosh \) case in a minute.)

But wait a minute, something is missing! We have 2 linearly independent solutions for \( Q(\phi) \), two linearly independent solutions for \( Z(z) \), but only one solution for \( R(r) \). Basic differential equation theory tells us we need 2 solutions for each of the 3 variables. Well it turns out, for non-integer \( \nu \), we have two Bessel functions: \( J_\nu \) and \( J_{-\nu} \). So the solution to Laplace’s equation is really

\[
R(r) = AJ_\nu(kr) + BJ_{-\nu}(kr)
\]

\[
\Phi(r, \phi, z) = (AJ_\nu(kr) + BJ_{-\nu}(kr)) \left( C \sin \nu \phi + D \cos \nu \phi \right) \left( E \sinh k z + F \cosh k z \right), \quad \nu = \text{non-integer}
\]

Finally, since Laplace’s equation is linear, the linear combination of any two solutions is also a solution:
\[ \Phi(r, \phi, z) = \sum_{\nu} \left( A_{\nu} J_{\nu}(kr) + B_{\nu} J_{\nu}(kr) \right) \left( C_{\nu} \sin \nu \phi + D_{\nu} \cos \nu \phi \right) \left( E_{\nu} \sinh \nu kz + F_{\nu} \cosh \nu kz \right), \]
\[\nu = \text{non-integer}\]

Note that each value of \( \nu \) has its own set of \( A, B, C, D, E, \) and \( F. \) Also, the values of \( \nu \) over which to sum are determined by the problem from the allowed range of \( \phi \) and the boundary conditions.

However, a very common case is \( \nu = \text{integer} \) (solution valid for \( \phi \) all around the axis). In that case, \( J_{\nu} \) is a multiple of \( J_{\nu} \), and the two solutions are not independent. So we do more math, and find a variant of the Bessel function which satisfies the equation, and is linearly independent. Such a function is a Neumann function (pronounced noy’-mon), also called a Bessel function of the second kind. Our solution is then:
\[ R(r) = AJ_{\nu}(kr) + BN_{\nu}(kr), \quad N_{\nu}(kr) \equiv \text{Neumann function}, \quad \nu = \text{integer}\]
\[ \Phi(r, \phi, z) = \sum_{\nu=0}^{\infty} \left( A_{\nu} J_{\nu}(kr) + B_{\nu} N_{\nu}(kr) \right) \left( C_{\nu} \sin \nu \phi + D_{\nu} \cos \nu \phi \right) \left( E_{\nu} \sinh \nu kz + F_{\nu} \cosh \nu kz \right) \]

An important property of the Bessel and Neumann functions is:

Bessel functions are regular at the origin, and Neumann functions blow up there.

This is analogous to the 2D polar case, where a domain that includes the origin must exclude the \( r^n \) solution, which blows up. In 3D, a cylinder that extends to \( r = 0 \) must exclude the Neumann function solution, which blows up at \( r = 0 \).

Case 2: For the case where \( Z(z) = \sin \) or \( \cos(kz) \), its eigenvalue is negative, and we have:
\[ \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = -\frac{\nu^2}{r^2} Q(\phi), \]
\[ \frac{\partial^2 Z}{\partial z^2} Z(z) = -k^2 Z(z) \]
\[ \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) R(r) = \left( +k^2 + \frac{\nu^2}{r^2} \right) R(r) \]

This is a different differential equation for \( R(r) \), and the Bessel functions don’t solve it. No problem; we just make up a new set of functions for this case, do the fancy math/power series solutions, and call them modified Bessel functions (not to be confused with Bessel functions of the 2nd or 3rd kinds). We use the same change of variable \( (w = kr) \), and call the solutions \( I_{\nu} \) and \( K_{\nu} \) [Jac ??]:
\[ \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) R(r) = \left( +k^2 + \frac{\nu^2}{r^2} \right) R(r) \quad \Rightarrow \quad R(r) = AI_{\nu}(kr) + BK_{\nu}(kr), \quad \nu \text{ integer} \]

As before, \( I_{\nu} \) is regular at the origin, and \( K_{\nu} \) blows up there. [For \( \nu \) non-integer, I assume we don’t use \( K_{\nu} \), and instead use \( I_{\nu} \) and \( I_{-\nu} \), as in the \( Z(z) = \sinh \) or \( \cosh \) case. Need reference??]

Three-D Laplace Solutions in Spherical Coordinates with Axial Symmetry

The Laplacian in spherical coordinates is
\[ \nabla^2 = \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) + \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) + \frac{1}{r^2 \sin^2 \theta} \frac{\partial^2}{\partial \phi^2}. \]

Note: \[ \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) = \frac{1}{r^2} \frac{\partial^2}{\partial r^2} (r \bullet) = \frac{2}{r} \frac{\partial}{\partial r} + \frac{\partial^2}{\partial r^2} \]

We start with a common special case of spherical coordinates: axial symmetry. In this case, the whole system is unchanged by any rotation about the z-axis. All of our results are thus independent of \( \phi \). We now seek a 2D solution \( \Phi(r, \theta) \), embedded in a 3D physical space. Applying the simplified Laplacian, and separation of variables, we seek something similar to the 2D polar case:

\[ \nabla^2 = \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) + \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) \quad \text{(with axial symmetry)} \]

\[ \Phi(r, \theta, \phi) = \Phi(r, \theta) = R(r)P(\theta) \]

As with the 2D polar case, the two operators must reduce to \( \pm k/r^2 \) for \( R(r) \) and \( P(\theta) \):

\[ \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) P(\theta) = -\frac{k}{r^2} P(\theta) \]

\[ \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) R(r) = \frac{k}{r^2} R(r) \]

The difference from the 2D polar case is that the differential operators for the two Laplacian terms are different. We don’t have any standard functions that provide the desired results, so (as with cylindrical coordinates), we invent them.

First, to find \( P(\theta) \): Fancy math \( \rightarrow \) **Legendre polynomials**, \( P_l(\cos \theta) \). These are a countably infinite set of polynomials, labeled with their (integer) order \( l \), which when applied to \( \cos \theta \), satisfy the above equation [Arf 11.37 p565]:

\[ \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) P_l(\cos \theta) = -\frac{l(l+1)}{r^2} P_l(\cos \theta) \]

The first few Legendre polynomials, and the corresponding \( P(\theta) \), are:

\[
\begin{align*}
P_0(x) &= 1, & P_1(x) &= x, & P_2(x) &= \frac{1}{2}(3x^2 - 1), & P_3(x) &= \frac{1}{2}(5x^3 - 3x) \\
P_4(x) &= \frac{1}{8}(35x^4 - 30x^2 + 3), & \\
P_1(\cos \theta) &= \cos \theta, & P_2(\cos \theta) &= \frac{1}{2}(3\cos^2 \theta - 1), & P_3(\cos \theta) &= \frac{1}{2}(5\cos^3 \theta - 3\cos \theta)
\end{align*}
\]

Now, to find \( R(r) \): Consider

\[ R(r) = Ar^l + Br^{-l-1} \quad \Rightarrow \quad \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) R(r) = \frac{1}{r^2} \frac{\partial}{\partial r} \left[ r^2 \left( Al^{l-1} - (l+1)Br^{-(l+2)} \right) \right] \]

\[ = \frac{1}{r^2} \frac{\partial}{\partial r} \left( Al^{l-1} - (l+1)Br^{-l} \right) = \frac{1}{r^2} \left( Al(l+1)r^l + (l+1)Br^{-(l+1)} \right) = \frac{l(l+1)}{r^2} R(r) \]

Finally, since Laplace’s equation is linear, the linear combination of any two solutions is also a solution:
\[ \Phi(r, \theta, \phi) = \Phi(r, \theta) = R(r)P(\theta) = \sum_{l=0}^{\infty} \left( A_l r^l + B_l r^{-l-1} \right) P_l(\cos \theta) \quad (\text{with axial symmetry}) \]

Note that each value of \( l \) has its own set of \( A \) and \( B \).

**Three-D Laplace Solutions in Spherical Coordinates without Axial Symmetry**

\[ \nabla^2 = \frac{1}{r^2} \frac{\partial}{\partial r} \left( r^2 \frac{\partial}{\partial r} \right) + \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) + \frac{1}{r^2 \sin^2 \theta} \frac{\partial^2}{\partial \phi^2} \]

\[ \Phi(r, \theta, \phi) = R(r)P(\theta)Q(\phi) \]

Without axial symmetry, we need a real function of \( \phi \). Since the prefactor of \( \partial^2 / \partial \phi^2 \) is \( 1/(r^2 \sin^2 \theta) \), which has no \( \phi \) in it, we simply need the eigenfunction of \( \partial^2 / \partial \phi^2 \), which is, as always, \( \sin \) or \( \cos(m\phi) \), with eigenvalue \(-m^2\):

\[ Q(\phi) = C \sin m\phi + D \cos m\phi \quad \Rightarrow \quad \frac{\partial^2}{\partial \phi^2} Q(\phi) = -m^2 Q(\phi). \]

The Legendre polynomials no longer work for \( \phi \), because now we need to satisfy

\[ \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) P(\theta) = \left( -\frac{l(l+1)}{r^2} + m^2 \right) P(\theta) \]

to cancel the new \(-m^2\) term from the \( \phi \) part above. For integer \( m \), more fancy math creates a new set of functions, the **associated Legendre functions**, \( P_l^m(\cos \theta) \), which satisfy the above equation for given \( l \) and \( m \):

\[ \frac{1}{r^2 \sin \theta} \frac{\partial}{\partial \theta} \left( \sin \theta \frac{\partial}{\partial \theta} \right) P_l^m(\cos \theta) = \left( -\frac{l(l+1)}{r^2} + m^2 \right) P_l^m(\cos \theta) \]

In other words, combining the \( \theta \) and \( \phi \) dependence, we see that (excluding \( r \) dependence):

\[ \nabla^2 P_l^m(\cos \theta)Q(\phi) = \left( -\frac{l(l+1)}{r^2} + m^2 \right) P_l^m(\cos \theta)Q(\phi) - m^2 P_l^m(\cos \theta)Q(\phi) = -\frac{l(l+1)}{r^2} P_l^m(\cos \theta)Q(\phi) \]

This is exactly the same equation we had for the axial symmetry case, so \( R(r) \) is not affected. Note that \( m \) is restricted to \(-l \leq m \leq l\). Also, when \( m = 0 \), the associated Legendre function is simply the original (non-associated) Legendre polynomial of order \( l \). [Associated Legendre functions are only polynomials for even \( l \); for odd \( l \), they have a square root in them.]

**The loss of axial symmetry introduced a \( \phi \) dependence, which was cancelled by new functions of \( \theta \). The radial function was not involved, and the radial function cannot “see” any dependence on \( \phi \).**

Now \( R(r) \): compare to the axially symmetric case (above): we added here a \(-m^2\) eigenvalue for \( \phi \), but cancelled all of it in the new \( P_l^m(\cos \theta) \). Therefore, the radial function doesn’t know anything about \( \phi \), and our axially symmetric \( R(r) \) still works as before.

\[ R(r) = A r^l + B r^{-l-1} \]

Finally, since Laplace’s equation is linear, the linear combination of any set of solutions is also a solution:
\[
\Phi(r, \theta, \phi) = R(r)P(\theta)Q(\phi) = \sum_{l=0}^{\infty} \left(A_l r^l + B_l r^{-l-1}\right) \sum_{m=-l}^{l} P_l^m(\cos \theta) \left(C_{lm} \sin m\phi + D_{lm} \cos m\phi\right)
\]

Note that each value of \(l\) has its own set of \(A_l\) and \(B_l\), and each \((l, m)\) pair has its own \(C_{lm}\) and \(D_{lm}\).

**Boundary Conditions Determine Solutions**

Given all our choices in the infinite set of solutions, how do we use boundary conditions to find the unique solution that they demand?

In rectangular coordinates of 2 or 3 dimensions, we have to choose between \(\sin/\cos\), \(\sinh/\cosh\), and \(e^x/e^{-x}\) for each coordinate. The boundary conditions determine the simplest form for Laplace solutions. It helps to consider the following facts about these functions:

<table>
<thead>
<tr>
<th>function</th>
<th>range</th>
<th>zeros at</th>
<th>derivative range</th>
<th>derivative range</th>
</tr>
</thead>
<tbody>
<tr>
<td>sin</td>
<td>-1 to 1</td>
<td>(n\pi)</td>
<td>cos</td>
<td>-1 to 1</td>
</tr>
<tr>
<td>cos</td>
<td>-1 to 1</td>
<td>((n+1/2)\pi)</td>
<td>-\sin</td>
<td>-1 to 1</td>
</tr>
<tr>
<td>sinh</td>
<td>-\infty to +\infty</td>
<td>0</td>
<td>\cosh</td>
<td>(\geq 1)</td>
</tr>
<tr>
<td>cosh</td>
<td>1 to (\infty)</td>
<td>none</td>
<td>\sin</td>
<td>-\infty to (\infty)</td>
</tr>
<tr>
<td>(\exp(+x))</td>
<td>&gt; 0 to (\infty)</td>
<td>none</td>
<td>(\exp(+x))</td>
<td>&gt; 0 to (\infty)</td>
</tr>
<tr>
<td>(\exp(-x))</td>
<td>&gt; 0 to (\infty)</td>
<td>none</td>
<td>-(\exp(-x))</td>
<td>-(\infty) to &lt; 0</td>
</tr>
</tbody>
</table>

(blue) \(\sinh(x)\) is odd, and has one zero. (green) \(\cosh(x)\) is even, and has no zeros. (red) \(e^{x/2}\) has no symmetry and no zeros.

Recall that in rectangular coordinates, at least one coordinate must be \(\sin/\cos\), and the other must be \(\sinh/\cosh\) or \(e^x/e^{-x}\). The following two cases often allow a simple choice for the coordinate which is \(\sin/\cos\):

1. Given \(\Phi(x=0, y) = 0\) and \(\Phi(x=b, y) = 0\).

   We try \(X(x) = A \sin(xc)\), because that is the only function above with zeros at \(x = 0\) and \(x = b\). Similarly for BCs in \(y\), and in 3D, for \(z\).

2. Given \(\partial/\partial x \Phi(x=0, y) = 0\) and \(\partial/\partial x \Phi(x=b, y) = 0\).
We try \( X(x) = A \cos(\pi x / b) \), because that is the only function above with zero derivatives at \( x = 0 \) and \( x = b \). Similarly for BCs in \( y \), and in 3D, for \( z \).

In 3D, we might find two coordinates which are both \( \sin/\cos \). Note, though, that other BCs can also be solved with \( \sin/\cos \), so the above choices are just two of many possible conditions.

Now having a good idea which coordinate(s) is/are \( \sin/\cos \), we use the other BCs with the table above to choose the simplest function for the other coordinate(s). These functions must be \( \sinh/\cosh \) or \( e^{kx}/e^{-x} \), because they have the positive eigenvalues needed to complement the negative ones from the \( \sin/\cos \) functions already identified in the solution. Some BCs allow us to write the solution as a single function, instead of a linear combination. For example:

(3) Given \( \Phi(x, y=0) = 0 \) and \( \Phi(x, y=b) = L \)

We try \( Y(y) = A \sinh(ky) \), because that is the only function above with zero at \( y=0 \) and non-zero at \( y=b \). Similarly for \( x \), and \( z \) in 3D.

(4) Given \( \Phi(x, y=0) = K \) and \( \Phi(x, y=b) = L \)

We try \( Y(y) = A \cosh(ky) \) or \( A \exp(ky) \), because those are the functions above non-zero at \( y=0 \) and at \( y=b \).

(5) Given \( \partial/\partial y \Phi(x, y=0) = 0 \) and \( \Phi(x, y=b) = L \)

We try \( Y(y) = A \cosh(ky) \), because that is the only function above with zero derivative at \( y=0 \), and non-zero value at \( y=b \).

(6) Given \( \partial/\partial y \Phi(x, y=0) = K \) and \( \partial/\partial y \Phi(x, y=b) = L \)

We try \( Y(y) = A \sinh(ky) \), because those are the functions above with non-zero derivatives at \( y=0 \), and at \( y=b \).

(7) Given symmetric BCs, such as \( \partial/\partial y \Phi(x, y = -a) = K \) and \( \partial/\partial y \Phi(x, y = +a) = K \)

We try \( Y(y) = A \cosh(ky) \), because it is the only symmetric function above with non-zero derivatives at \( y = \pm a \).

(8) Given anti-symmetric BCs, such as \( \partial/\partial y \Phi(x, y = -a) = -K \) and \( \partial/\partial y \Phi(x, y = +a) = +K \)

We try \( Y(y) = A \sinh(ky) \), because it is the only anti-symmetric function above with non-zero derivatives at \( y = \pm a \).

If the BCs do not allow any of the above simple solutions, then the solution must be a linear combination of these functions, with coefficients that must be determined by solving simultaneous equations, in the usual way.

Respecting Orthogonality

TBS: orthogonality wrt to linear operators. Orthogonality of interest of Bessel functions.

Propagation In a Vacuum

Maxwell’s equations imply that all vacuum EM fields, including inside waveguides, satisfy the wave equations (in Gaussian units) [Jac sec. 8.2]:
\[ \nabla^2 \mathbf{E} = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} \mathbf{E}, \quad \] i.e., each component of \( \mathbf{E} \) satisfies the wave equation:
\[ \nabla^2 E_x = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} E_x, \quad \nabla^2 E_y = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} E_y, \quad \text{and} \quad \nabla^2 E_z = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} E_z. \]

Also for \( \mathbf{B} \):
\[ \nabla^2 \mathbf{B} = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} \mathbf{B} \]
\[ \Rightarrow \quad \nabla^2 B_x = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} B_x, \quad \nabla^2 B_y = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} B_y, \quad \text{and} \quad \nabla^2 B_z = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} B_z. \]

---

In free space (no waveguide), the \( \mathbf{E} \) & \( \mathbf{B} \) fields are in phase.

Many people think that in free space, \( \mathbf{E} \) & \( \mathbf{B} \) are out of phase, and imagine the energy to be shifted back and forth between the \( \mathbf{E} \) & \( \mathbf{B} \) fields, but that is not true.

Absorbing an EM wave is like being hit with a stream of warm beans. The energy density at the \( \mathbf{E} \) & \( \mathbf{B} \) nodes (zeros) is zero. The energy density is maximum at the peaks (anti-nodes).

---

Linearly polarized EM wave in free space, propagating in the \( z \) direction. On the right, only the \( \mathbf{E} \) and \( \mathbf{B} \) fields are shown, for simplicity. On the left, the \( \mathbf{A} \) field is included. \( \mathbf{A} \) is in the plane of \( \mathbf{E} \), but leads \( \mathbf{E} \) in time and space.

In the radiation gauge, the vector potential \( \mathbf{A} \) for a linearly polarized EM wave is also a linearly polarized transverse wave, in the plane of the \( \mathbf{E} \) field. \( \mathbf{A} \) leads \( \mathbf{E} \) in time (\( \mathbf{A} \) peaks 90° before \( \mathbf{E} \)), and lags in space (\( \mathbf{A} \) peaks 90° after \( \mathbf{E} \)). Consider the point \( w \) in space in the diagram above:

\[ \mathbf{A} = A_y \hat{\mathbf{y}}, \quad \mathbf{A}(w) = 0, \quad \frac{\partial A_y(w)}{\partial t} < 0, \quad \frac{\partial A_y(w)}{\partial z} > 0. \]

\( \mathbf{E} \) is maximum at \( w \), where \( \mathbf{A} \) is 0 and decreasing in time:

\[ \mathbf{E} = -\frac{1}{c} \frac{\partial \mathbf{A}}{\partial t} \]

At the same point, \( \mathbf{B} \) is (negative, in this case) maximum at \( w \), where \( \mathbf{A} \) is 0 and increasing in space:
\[
\mathbf{B} = \nabla \times \mathbf{A} = \left( \frac{\partial A_z}{\partial y} - \frac{\partial A_y}{\partial z} \right) \hat{x} = -\frac{\partial A_y}{\partial z} \hat{x}
\]

However, the \( \mathbf{A} \) field decreases in magnitude with increasing frequency, for the same power density, because \( \mathbf{E} \) & \( \mathbf{B} \) derive from \( \mathbf{A} \) with derivative operators, which introduce factors of \( \omega \) for \( \mathbf{E} \), or \( k \) for \( \mathbf{B} \):

**E vs. A:**

- Low Frequency, \( \mathbf{A} \) larger
- Low Frequency, \( \mathbf{A} \) smaller

![Same magnitude E field and its associated A field at two different frequencies.](image)

**Waveguides**

We should first note that in the context of waveguides, the word “mode” is somewhat abused: in classical mechanics, and in E&M cavities, a “mode” implies a specific frequency of oscillation. In waveguides, the “mode” really refers to the eigenvalue of the waveguide transverse eigenfunction/eigenvalue equation (described below), and covers a continuum of temporal frequencies, \( \omega \), and spatial frequencies (wave numbers), \( k \). One could say the waveguide “mode” refers only to the transverse mode, but not the whole of the EM wave(s) in the waveguide.

This section relies heavily on the methods developed in “Solving Laplace’s Equation” and “Propagation in a Vacuum,” above. You should read those sections first. Those sections concern electrostatics and magnetostatics. In the topic of waveguides, we extend those principles to electrodynamics, where the fields are time-dependent. We therefore use the full, time-dependent Maxwell equations, including time derivatives.

Waveguides are concerned with the propagation of waves down a fixed (but arbitrary) cross sectioned cylinder. The most common cross sections are rectangles, circles, and fractions of a circle. Since we are concerned with wave propagation, prior experience strongly suggests decomposition of an arbitrary wave into sinusoidal components. [More rigorously, sinusoids compose a basis set of functions for the construction of any arbitrary function, and even better, are the eigenfunctions of linear integro-differential equations (such as the wave equations resulting from Maxwell’s equations). This means we can study a single sinusoid of arbitrary frequency, and know that an arbitrary function can be made from a linear superposition of these sinusoids, and the sinusoids will not interact with each other. This is the standard method of eigenfunction decomposition and analysis, familiar from Fourier analysis and quantum mechanics.]

Under what conditions can a sinusoidal wave propagate down the waveguide, but remain otherwise unchanged? We assume an ideal conductor for the boundary of the waveguide, so there is no loss as the wave propagates. (Loss is a small correction to the lossless case [Jac p353m]). The choice of propagating sinusoids leads immediately to the use of phasors (complex numbers) for all components of \( \mathbf{E} \) and \( \mathbf{B} \) fields; WLOG, we assume propagation in the +\( z \) direction. Then all 3 components of both the \( \mathbf{E} \) and \( \mathbf{B} \) fields vary sinusoidally in time and \( z \)-direction:

\[
\mathbf{E}(x, y, z, t) = \text{Re}\left\{ \mathbf{E}(x, y)e^{ikz-\omega t} \right\}, \quad \mathbf{B}(x, y, z, t) = \text{Re}\left\{ \mathbf{B}(x, y)e^{ikz-\omega t} \right\} \quad \text{[Jac 8.18 p357]}
\]
Note that $\omega$ is always real, and for lossless propagation, $k$ must be real. So far, these are still fully 3-dimensional vector fields with components in all 3 directions. All the components propagate uniformly in the $z$ direction. Because we have now defined the variation of the fields in $z$ and $t$, we can find the time and $z$ derivatives, and partially evaluate the Laplacian operator in the wave equations above:

$$\frac{\partial^2 E}{\partial z^2} = -k^2 E, \quad \frac{\partial^2 E}{\partial t^2} = -\omega^2 E, \quad \frac{\partial^2 B}{\partial z^2} = -k^2 B, \quad \frac{\partial^2 B}{\partial t^2} = -\omega^2 B.$$ 

Since we know the variation in $z$, it becomes useful to separate the Laplacian operator into its "$z$-part" and its "transverse part" (i.e., its $x$-$y$ part), so we define:

$$\nabla^2 = \nabla_t^2 + \frac{\partial^2}{\partial z^2} \quad \Rightarrow \quad \nabla_t^2 = \nabla^2 - \frac{\partial^2}{\partial z^2} = \frac{\partial^2}{\partial t^2} + \frac{\partial^2}{\partial z^2} = \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} \quad \text{[Jac 8.20 p357]}. $$

Note the subscript "t" means "transverse" here, not "tangential." Also, we may use polar coordinates $(r, \phi)$ or $(x, y)$ coordinates for the transverse part (or any other 2-D coordinates for that matter), whichever is easier. With these definitions, keeping in mind that all vector components are complex-valued phasors, the wave equations above become:

$$\left( \nabla_t^2 + \frac{\partial^2}{\partial z^2} \right) E = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} E \quad \Rightarrow \quad \left( \nabla_t^2 - k^2 \right) E = -\frac{\omega^2}{c^2} E, \quad \nabla_t^2 E = \left( k^2 - \frac{\omega^2}{c^2} \right) E,$$

or in components:

$$\nabla_t^2 E_x = \left( k^2 - \frac{\omega^2}{c^2} \right) E_x \quad \nabla_t^2 E_y = \left( k^2 - \frac{\omega^2}{c^2} \right) E_y,$$

and

$$\nabla_t^2 E_z = \left( k^2 - \frac{\omega^2}{c^2} \right) E_z.$$

Also for $B$:

$$\left( \nabla_t^2 + \frac{\partial^2}{\partial z^2} \right) B = \frac{1}{c^2} \frac{\partial^2}{\partial t^2} B \quad \Rightarrow \quad \nabla_t^2 B = \left( k^2 - \frac{\omega^2}{c^2} \right) B,$$

i.e.,

$$\nabla_t^2 B_x = \left( k^2 - \frac{\omega^2}{c^2} \right) B_x \quad \nabla_t^2 B_y = \left( k^2 - \frac{\omega^2}{c^2} \right) B_y,$$

and

$$\nabla_t^2 B_z = \left( k^2 - \frac{\omega^2}{c^2} \right) B_z.$$

Thus we see that the conditions for propagation down a waveguide are that all 6 E & M components, $E_x, E_y, E_z, B_x, B_y,$ and $B_z$, must satisfy the same 2D eigenfunction/eigenvalue equation in the transverse plane. Note, however, that some of the 6 components may be everywhere zero, in which case those components identically satisfy any (linear) eigenfunction/eigenvalue equation. Each scalar field eigenfunction has an eigenvalue, which determines the dispersion relation relating $k$ and $\omega$ for that eigenfunction. Further, to achieve propagation of such a mode, $k$ and $\omega$ must be the same for all 6 E & M components. Therefore the eigenvalues (and thus dispersion relation) must be the same for all non-zero E- and B-field components of a single frequency. It is common to define the eigenvalue for a given waveguide frequency, and therefore the dispersion relation, as:

$$\text{eigenvalue} = -\gamma^2 = k^2 - \frac{\omega^2}{c^2}, \quad \gamma^2 = \frac{\omega^2}{c^2} - k^2 \quad \Rightarrow \quad \omega(k) = c \sqrt{k^2 + \gamma^2}.$$
Dispersion relation for EM propagation in a waveguide.

Note that for hollow cross-sectional waveguides, the eigenvalues are always negative (why??), so $\gamma^2 > 0$. For propagation, $k$ must be real $> 0$. Thus, a given propagation mode has a minimum frequency for propagation of

$$\omega_{\text{min}} = c\gamma .$$

In free space, and with plane or spherical waves, $\gamma = 0$, the dispersion relation is linear: $\omega = ck$, and there is no frequency dispersion. Note that as $\omega \to \infty$, the dispersion relation for all waveguide modes is asymptotic to $\omega = ck$, which is the free space dispersion relation (of no actual dispersion).

Now that we have the basic equations of operation of a waveguide, we can consider the boundary conditions that exist in a waveguide.

**Boundary Conditions and Propagation**

The walls of a waveguide are good conductors, which imposes boundary conditions that further constrain the propagation of EM waves down the guide, compared to free-space propagation. In particular, a good conductor has surface charges and currents that eliminate all $E$ (& $D$) and $B$ (& $H$) fields inside it (except for a small skin depth) [Jac p358]. Also, as always, the $E_{\text{parallel}} (= E_z)$ and $B_n (= B_{\text{transverse}})$ are continuous at the boundaries. Therefore, at the boundaries,

$$E_z = 0, \quad B \cdot \hat{n} = 0. \quad (NB: B_t = B_x \hat{x} + B_y \hat{y}).$$

Note the subscript “t” still means “transverse” here, not “tangential.”

TBS.

**Phase and Group Velocity In a Waveguide**

First, consider a plane wave propagating in free space at an angle to the $z$-axis (Figure 3.11, left).

**Figure 3.11** (Left) Phase velocity in free space. (Right) Phase and group velocity in a waveguide.
In the direction of propagation (along the red arrow), the phase velocity is just $c$, the speed of light. But suppose we set up electrometers to measure the electric field as it passes down the $z$-axis. The time between wave crests is the same, but the *distance* between wave crests along the $z$-axis is greater. So the wave crests are moving *faster than light* along the $z$-axis; this is required by the geometry. Hence, in the $z$-direction, $v_{\text{phase}} > c$.

The propagation down a waveguide can be written as a superposition of two constituent waves traveling at an angle to the waveguide axis (Figure 3.11, right, blue and red). The constituent waves reflect off the walls of the waveguide, and superpose to produce a net propagation in the $z$ direction. The phase velocity of both the blue and red constituents is the same as in the left diagram:

$$v_{\text{phase}} = c / \cos \theta > c .$$

However, the net propagation speed in the $z$ direction is just the $z$ component of the red (or blue) constituent:

$$v_{\text{group}} = c(\cos \theta) < c .$$

Therefore, in a waveguide, the phase velocity is always $> c$, and the group velocity is always $< c$.

Note that the constituents are *not* simply two uniform plane waves. The boundary conditions must be met on each constituent wave: namely, that $E_r = 0$ at the walls, and that $B_r = 0$ (the normal component of the B-field). So the constituent waves also taper to zero at the edges, just like the total wave does.??

The angle $\theta$ can be determined from the $z$-component of the spatial frequency (the wave-vector), $k_z$. We find $k_z$ from solving Maxwell’s equations, and $\theta$ from:

$$k_z = -\frac{\omega}{c \cos \theta} .$$

**Multiple modes of one frequency:** It is generally the case that a single frequency can propagate in multiple modes. The modes typically have different group velocities, though, so they are dispersed in time and space as they propagate. This is usually undesirable. To avoid this, the transmitter must excite only a single desired mode. Methods of doing this are beyond our scope.

**Cylindrical Hollow Waveguides**

![Cylindrical waveguide](image)

**Figure 3.12** Cylindrical waveguide.

Cylindrical waveguides are common: they are just metal pipes, possibly gently bent to guide the waves along a path. The propagation modes are different than those of rectangular waveguides, but are derived from the same principles. In cylindrical coordinates, just as in rectangular, we generalize Laplace’s equation (3.1) to time-dependent fields by including the time-dependent factor $e^{-i\omega t}$. This gives the wave equation:

$$\nabla^2 B_z = -\frac{\omega^2}{c^2} E_z , \quad \nabla^2 = \frac{1}{r} \frac{\partial}{\partial r} \left( r \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} + \frac{\partial^2}{\partial z^2} \quad \Rightarrow \quad B_z(r, \phi, z) = R(r)Q(\phi)Z(z)e^{-i\omega t} \quad \text{(and similar for } E_z).$$

As before, we seek propagation modes down the pipe, so:
\[ Z(z) = e^{ik_z z}, \quad \text{and} \quad \frac{\partial^2}{\partial z^2} Z(z) = -k_z^2 Z(z) \quad \text{where} \quad k_z \equiv \text{wave number}. \]

**Solutions:** As when solving Laplace’s equation, the azimuthal function must satisfy:

\[
\frac{1}{r^2} \frac{\partial^2}{\partial \phi^2} Q(\phi) = -\frac{k_{\phi}^2}{r^2} Q(\phi) \Rightarrow \frac{\partial^2}{\partial \phi^2} Q(\phi) = -k_{\phi}^2 Q(\phi), \quad Q(\phi) = e^{ik_{\phi} \phi}.
\]

It is conventional to call \( k_{\phi} \) ‘\( m \)’: \( m \equiv k_{\phi}, \quad Q(\phi) = e^{im\phi}. \)

Then \( R(r) \) must satisfy:

\[
\frac{1}{r} \frac{\partial}{\partial r} \left(r \frac{\partial}{\partial r} \right) R(r) = \left( -k_r^2 + \frac{m^2}{r^2} \right) R(r), \quad \text{or} \quad \left[ \frac{1}{r} \frac{\partial}{\partial r} \left(r \frac{\partial}{\partial r} \right) + k_r^2 - \frac{m^2}{r^2} \right] R(r) = 0.
\]

Notice that the second term on the right cancels the azimuthal term from the \( \phi \) contribution. As with Laplace’s equation, this equation is singular at \( r = 0 \), which is often a physically valid radius. The singularity is easily removed by defining \( R(r = 0) \) as the limit:

\[ R(0) \equiv \lim_{r \to 0} R(r). \]

The differential equation is the famous Bessel equation, solved by Bessel functions of the first kind, \( J_m(x) \):

\[ R(r) = J_m(k_r r). \]

**Quantization:** In the rectangular case, \( k_x \) and \( k_y \) were quantized by the boundary conditions of the guide. In the cylindrical case, \( k_z \) and \( m \) are quantized. With the cancellation of the azimuthal terms, \( k_r^2 \) acts like \( (k_x^2 + k_y^2) \), and puts a lower limit \( \omega_{\text{min}} \) on possible solutions to Maxwell’s equations in the guide. The constraint is:

\[ -k_r^2 - \beta^2 = -\frac{\omega^2}{c^2} \quad \text{or} \quad \beta^2 = \frac{\omega^2}{c^2} - k_r^2 > 0. \]

The azimuthal parameter \( m \) must make the wave-function single-valued as \( \phi \) goes across \( 2\pi \). Therefore, \( m \) must be an integer. (This is exactly the same quantization and azimuthal function as cylindrical and spherical systems in quantum mechanics.) **In our case, we can take** \( m \geq 0 \), so:

\[ m = 0, 1, 2, \ldots \]

In preparation for considering the radial modes, we recall the form of Bessel functions of first kind (Figure 3.13). They oscillate, with an infinite number of zeros, and extrema.
Figure 3.13  Some Bessel functions of the first kind.

The radial function $R(r)$ is quantized by the boundary condition at the walls of the pipe. For TE modes, $B_t$ must be stationary at the edge $r = a$, so:

$$\frac{\partial B_z}{\partial r} \propto \frac{dR(a)}{dr} = 0 \Rightarrow J_m'(k_r a) = 0, \quad k_r = \frac{x_m}{a}, \quad x_m \equiv j^{th} \text{ zero of } J_m, \quad j = 1, 2, ...$$

$$B_z(t, r, \phi, z) = J_m(k_r r)e^{i \phi} e^{i(k_z - \omega t)}, \quad E_z(t, r, \phi, z) = 0.$$

These solutions are the $TE_{mj}$ modes.

For TM modes, $E_z = 0$, so $E_t = 0$, and:

$$E_z(z) \propto R(a) = 0 \Rightarrow J_m(k_r a) = 0, \quad k_r = \frac{x_m}{a}, \quad x_m \equiv j^{th} \text{ zero of } J_m, \quad j = 1, 2, ...$$

$$E_z(t, r, \phi, z) = J_m(k_r r)e^{i \phi} e^{i(k_z - \omega t)}, \quad B_z(t, r, \phi, z) = 0.$$

These solutions are the $TM_{mj}$ modes.

Order of the modes: Within the TE cases, and within the TM cases, except for the $m = 0$ case, $k_r$ increases as $m$ increases, i.e. $k_r$ increases with the azimuthal mode, so the cutoff frequency increases with both $m$ and $j$, similar to the rectangular waveguide. You can see from Figure 3.13 that $J_0$ is special, because mathematically, its first stationary point is $x = 0$, but that is not a physical mode. Therefore, the fastest mode (aka “dominant mode”, with smallest $k_r$) is $TE_{11}$. However, the TE modes interleave with TM modes, so the sequence of modes with increasing cutoff frequency $\omega_{\text{min}}$ is somewhat awkward:

<table>
<thead>
<tr>
<th>$m$</th>
<th>$TE (x'_{mj})$</th>
<th>$TM (x_{mj})$</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>3.8317</td>
<td>2.4048</td>
</tr>
<tr>
<td>1</td>
<td>1.8412</td>
<td>3.8317</td>
</tr>
<tr>
<td>2</td>
<td>3.0542</td>
<td>5.1356</td>
</tr>
<tr>
<td>3</td>
<td>4.2012</td>
<td>6.3802</td>
</tr>
</tbody>
</table>

**Multipoles: Dipoles and Quadrupoles**

Suppose we have a small blob of charges, and we look at it from a distance. We are at the origin:
We wish to know the potential we feel (here at the origin) from this blob of charges. From that, we can compute the E-field. We suppose that our observation distance is large compared to the size of the blob, as shown, so we can use the far field approximation. This is a valuable and widely used approximation. We assume the blob consists of \( n \) point charges, with total charge

\[
Q_{\text{total}} = \sum_{i=1}^{n} Q_i.
\]

To zeroth order, the potential is simply

\[
\Phi = \frac{Q_{\text{total}}}{r}, \quad \text{where} \quad r = |r|.
\]

If all the point charges were exactly at \( r \), this would be the exact potential. Now consider the effect of moving one of the point charges, \( Q_i \), from the point \( r \) by some small amount, \( r' \). The potential at the origin due to that charge changes by

\[
\Delta \Phi_{Q_i} = \frac{Q_i}{|r + r'|} - \frac{Q_i}{r} = Q_i \left( \frac{1}{|r + r'|} - \frac{1}{r} \right) = Q_i \left( \frac{1}{r} \right) \left( \frac{1}{r'} \right).
\]

That is, the change in potential (seen by us at the origin) equals the charge times the change in \( 1/r \). The function \( 1/r \) is a scalar field: it assigns a number to every point in space. Therefore, we can use its gradient to approximate small changes in \( 1/r \):

\[
\Delta \frac{1}{r} \approx \nabla \left( \frac{1}{r} \right) \cdot r' \quad \Rightarrow \quad \Delta \Phi_{Q_i} \approx Q_i \nabla \left( \frac{1}{r} \right) \cdot r'.
\]

We can find the gradient direction with a simple geometric argument: the gradient points in the direction of greatest rate of increase.

The greatest rate of increase of \( 1/r \) is opposite to \( r \) itself; \( r \) increases fastest along itself. Therefore, \( 1/r \) decreases fastest along \( r \), and \( 1/r \) increases fastest opposite to \( r \): The magnitude of the gradient then follows from a simple derivative:

\[
\frac{d}{dr} \frac{1}{r^2} = -\frac{1}{r^2} \quad \Rightarrow \quad \nabla \left( \frac{1}{r} \right) = -\frac{1}{r^2} \hat{r}.
\]
We use this to approximate the change in potential at the origin due to moving \( Q_i \) a small distance \( r' \):

\[
\Delta \Phi_{Q_i} = Q_i \nabla \left( \frac{1}{r} \right) \cdot r' = Q_i \frac{-1}{r'^2} \hat{r} \cdot r'.
\]

Now let’s move each of the charges by its own small displacement \( r'_i \). By the principle of superposition, we just add up the changes due to moving each charge:

\[
\Delta \Phi_{\text{total}} = \sum_{i=1}^{n} Q_i \frac{-1}{r'^2} \hat{r} \cdot r'_i = \frac{-1}{r'^2} \hat{r} \cdot \left( \sum_{i=1}^{n} Q_i r'_i \right) = \frac{-1}{r'^2} \hat{r} \cdot p, \quad \text{where } p = \sum_{i=1}^{n} Q_i r'_i.
\]

Note that \( p \) is independent of where the blob is located, i.e. independent of the relative position of observer and blob. It is solely a function of the blob, i.e. the distribution of charges. \( p \) is called the dipole moment of the charge distribution. We see that, to first order, the potential due to a net-neutral blob (I mean, charge distribution) falls off as \( 1/r^2 \), and is zero along a line through the blob perpendicular to \( p \).

Note that often we consider the blob to be at the origin, and compute the field at some position \( r \). This simply reverses the sign of \( r \) above, to yield the more common formula:

\[
\Delta \Phi_{\text{total}} = \frac{1}{r'^2} \hat{r} \cdot p \Rightarrow \Phi \approx \frac{Q_{\text{total}}}{r} + \frac{1}{r'^2} \hat{r} \cdot p + \ldots.
\]

A subtle point: the dipole moment of the blob may depend on the reference point within the blob from which the \( r' \) are measured. However, if the total charge is zero, the dipole moment is then independent of the reference point within the blob. Why??

Note that the dipole moment does not completely characterize the charge distribution, nor is it necessarily the most significant component of such a characterization. It is simply the \( 1/r^2 \) component of the potential due to the charge distribution in the far field approximation.

**Quadrupoles**

The quadrupole moment extends the above approximation to \( 2^\text{nd} \) order. In short, we could say, “Just extend the dipole to the 3-variable \( 2^\text{nd} \) order Taylor expansion of \( 1/r^2 \),” and claim we’re done. But that’s too arcane, so let’s see what it really means.

First, let’s recall the meaning of the Taylor expansion of a simple function, \( f(x) \), about \( x_0 \). To first order, we approximate the first derivative as constant. Hence

\[
\Delta f \approx f'(x_0) \Delta x \Rightarrow f(x + \Delta x) \approx f(x_0) + f'(x_0) \Delta x.
\]

To \( 2^\text{nd} \) order, we approximate the \( 2^\text{nd} \) derivative as constant, and therefore the first derivative changes linearly with \( x \):

\[
\Delta f' \approx f''(x_0) \Delta x \Rightarrow f'(x + \Delta x) \approx f'(x_0) + f''(x_0) \Delta x.
\]

We can then find the average first derivative \( f' \) over the interval \( \Delta x \):

\[
f'_{\text{avg}} = f'(x_0) + \frac{1}{2} \Delta f' = f'(x_0) + \frac{1}{2} f''(x_0) \Delta x.
\]

This average \( f' \) incorporates \( 2^\text{nd} \) order effects. Now use this average \( f' \) in the first order estimate for \( f(x + \Delta x) \):

\[
\Delta f \approx f'_{\text{avg}}(x_0) \Delta x = \left( f'(x_0) + \frac{1}{2} f''(x_0) \Delta x \right) \Delta x = f'(x_0) \Delta x + \frac{1}{2} f''(x_0) (\Delta x)^2
\]

\[
\Rightarrow f(x + \Delta x) \approx f(x_0) + f'(x_0) \Delta x + \frac{1}{2} f''(x_0) (\Delta x)^2.
\]

Dipole moments are the \( 1^\text{st} \) order Taylor coefficients in the expansion of potential due to a blob of charges. Quadrupole moments are simply the \( 2^\text{nd} \) order Taylor coefficients in the same expansion:
\[ \Phi(\mathbf{r}) = \frac{Q_{\text{total}}}{r} + \frac{\mathbf{p} \cdot \hat{\mathbf{r}}}{r^2} + \frac{\text{quadrupole term}}{r^4} + \ldots \]

We follow a procedure similar to the 2\textsuperscript{nd} order Taylor expansion of \( f(x) \) above, but with two additional complications: (1) our function \( 1/r \) is a function of 3-D space, so its first derivative (gradient) is a vector, not a number; and (2) the 2\textsuperscript{nd} derivative of \( 1/r \) is the gradient of a gradient, which is the gradient of a vector, which is a rank-2 tensor. We take these two issues in turn. As is traditional, we’ll stick to Cartesian coordinates.

[Aside: purists might object to the following explanation, but our goal here is to describe electromagnetics, not differential geometry.]

First, the gradient of a scalar field at a point is a set of 3 numbers (a vector), which tell us how the scalar field varies as we move in the \( x \), \( y \), and \( z \) directions:

\[ \nabla f = \left( \frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial f}{\partial z} \right). \]

We usually write the gradient as a vector, but for now, you can think of it as just a set of 3 numbers. We use the 3 numbers to approximate changes in \( f \) given small changes in position, \( dx \), \( dy \), \( dz \):

\[ \Delta f \approx \frac{\partial f}{\partial x} \Delta x, \quad \Delta f \approx \frac{\partial f}{\partial y} \Delta y, \quad \Delta f \approx \frac{\partial f}{\partial z} \Delta z. \]

Further, for differentially small changes \( dx \), \( dy \), \( dz \), the \( \Delta f \) are independent and can be summed, to get the familiar total derivative rule:

\[ \Delta f \approx \frac{\partial f}{\partial x} \Delta x + \frac{\partial f}{\partial y} \Delta y + \frac{\partial f}{\partial z} \Delta z. \]

This is what we used to compute the dipole moment. It is a first order approximation, because we take the gradient to be constant over the intervals \( dx \), \( dy \), \( dz \). To make the 2\textsuperscript{nd} order approximation, we take the 2\textsuperscript{nd} derivative to be constant, thus making the gradient (first derivative) vary linearly over the intervals \( dx \), \( dy \), \( dz \).

So now we ask, what is the gradient of a vector field, \( \mathbf{g}(\mathbf{r}) \)? At a given point, it is simply a set of 3 vectors that tell us how the vector field varies as we move in the \( x \), \( y \), and \( z \) directions:

\[ \nabla \mathbf{g} = \left( \frac{\partial \mathbf{g}}{\partial x}, \frac{\partial \mathbf{g}}{\partial y}, \frac{\partial \mathbf{g}}{\partial z} \right) \]

\[ \Delta \mathbf{g} \approx \left( \frac{\partial \mathbf{g}}{\partial x} \right) \Delta x + \left( \frac{\partial \mathbf{g}}{\partial y} \right) \Delta y. \]

\[ \Delta \mathbf{g} \approx \left( \frac{\partial \mathbf{g}}{\partial x} \right) \Delta x + \left( \frac{\partial \mathbf{g}}{\partial y} \right) \Delta y. \]

[Figure 3.15 Variation of a 2D vector field over small intervals.]

Note that all 3 component of \( \mathbf{g} \) may vary even if we move only in, say, \( x \). We can approximate the change in the vector \( \mathbf{g} \) over small displacements, just like we did for the scalar field:
\[ \Delta \mathbf{g} \approx \frac{\partial \mathbf{g}}{\partial x} \, dx, \quad \Delta \mathbf{g} \approx \frac{\partial \mathbf{g}}{\partial y} \, dy, \quad \Delta \mathbf{g} \approx \frac{\partial \mathbf{g}}{\partial z} \, dz, \quad \text{and} \quad \Delta \mathbf{g} \approx \frac{\partial \mathbf{g}}{\partial x} \, dx + \frac{\partial \mathbf{g}}{\partial y} \, dy + \frac{\partial \mathbf{g}}{\partial z} \, dz. \]

\( \nabla \mathbf{g} \) is a set of 3 vectors, each with 3 components. We write \( \mathbf{g} \) as a 3 x 3 matrix of 3 column vectors:

\[
\nabla \mathbf{g} = \left( \frac{\partial \mathbf{g}}{\partial x}, \frac{\partial \mathbf{g}}{\partial y}, \frac{\partial \mathbf{g}}{\partial z} \right) = \begin{bmatrix} g_{xx} & g_{yx} & g_{zx} \\ g_{xy} & g_{yy} & g_{zy} \\ g_{xz} & g_{yz} & g_{zz} \end{bmatrix}, \quad \text{where} \quad \frac{\partial \mathbf{g}}{\partial x} = \begin{bmatrix} g_{xx} \\ g_{xy} \\ g_{xz} \end{bmatrix}, \quad \frac{\partial \mathbf{g}}{\partial y} = \begin{bmatrix} g_{yx} \\ g_{yy} \\ g_{yz} \end{bmatrix}, \quad \frac{\partial \mathbf{g}}{\partial z} = \begin{bmatrix} g_{zx} \\ g_{zy} \\ g_{zz} \end{bmatrix}.
\]

This is a rank-2 tensor. It can operate on a displacement vector to approximate the change in a vector field over that displacement

\[
\Delta \mathbf{g} \approx \frac{\partial \mathbf{g}}{\partial x} \, dx + \frac{\partial \mathbf{g}}{\partial y} \, dy + \frac{\partial \mathbf{g}}{\partial z} \, dz \equiv \nabla \mathbf{g} (dx) = \begin{bmatrix} g_{xx} & g_{yx} & g_{zx} \\ g_{xy} & g_{yy} & g_{zy} \\ g_{xz} & g_{yz} & g_{zz} \end{bmatrix} (dx).
\]

Note that a tensor is linear in its vector argument: \( \nabla \mathbf{g} (k \, dx_1 + dx_2) = k \nabla \mathbf{g} (dx_1) + \nabla \mathbf{g} (dx_2) \). [This implies that if we transform coordinates (e.g., rotation), the columns and then rows of the tensor will transform like vectors; but we don’t really care about that right now.]

Now let’s compute the gradient of the gradient of a scalar function \( f(x, y, z) \):

\[
\nabla f = \left( \frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial f}{\partial z} \right) = \begin{bmatrix} \frac{\partial^2 f}{\partial x^2} \\ \frac{\partial^2 f}{\partial y^2} \\ \frac{\partial^2 f}{\partial z^2} \end{bmatrix}.
\]

\[

abla \nabla f = \begin{bmatrix} \frac{\partial^2 f}{\partial x^2} & \frac{\partial^2 f}{\partial x \partial y} & \frac{\partial^2 f}{\partial x \partial z} \\ \frac{\partial^2 f}{\partial y \partial x} & \frac{\partial^2 f}{\partial y^2} & \frac{\partial^2 f}{\partial y \partial z} \\ \frac{\partial^2 f}{\partial z \partial x} & \frac{\partial^2 f}{\partial z \partial y} & \frac{\partial^2 f}{\partial z^2} \end{bmatrix}.
\]

It’s just the matrix of all second derivatives. [It is called the Hessian matrix.]

Back to quadrupoles: We are the observer at the origin. We want the 2nd order change in potential seen by us due to displacing a single charge \( Q \) from the blob by a small amount \( \mathbf{r}' = (x', y', z') \). We follow the same procedure as above for the 2nd order Taylor expansion of \( f(x) \). First, we compute the gradient of the gradient of \( 1/r \):
\[ \nabla \nabla \left( \frac{1}{r} \right) = \nabla \left[ -\frac{1}{r^3} \hat{r} \right] = \nabla \left[ -\frac{1}{r^3} \right] = \nabla \left[ \left( x^2 + y^2 + z^2 \right)^{-3/2} (\hat{x} \hat{y} + \hat{z}) \right]. \]

Use \( \nabla \Phi \equiv \left( \frac{\partial \Phi}{\partial x}, \frac{\partial \Phi}{\partial y}, \frac{\partial \Phi}{\partial z} \right) \)

\[ \frac{\partial}{\partial x} \left( x^2 + y^2 + z^2 \right)^{-3/2} (\hat{x} \hat{y} + \hat{z}) \]

\[ = \frac{3}{2} \left( x^2 + y^2 + z^2 \right)^{-5/2} \left( \hat{x} x + \hat{y} y + \hat{z} z \right) \]

\[ = \left[ \frac{3x^2}{r^5} - \frac{1}{r^3} \right] \hat{x} + \frac{3xy}{r^5} \hat{y} + \frac{3xz}{r^5} \hat{z} \]

Similarly: \[ \frac{\partial}{\partial y} \left[ \frac{1}{r^3} \right] = \frac{3xy}{r^5} \hat{x} + \frac{3y^2}{r^5} \hat{y} + \frac{3yz}{r^5} \hat{z}, \quad \text{and} \quad \frac{\partial}{\partial z} \left[ \frac{1}{r^3} \right] = \frac{3zx}{r^5} \hat{x} + \frac{3zy}{r^5} \hat{y} + \frac{3z^2}{r^5} - \frac{1}{r^3} \hat{z} \]

\[ \nabla \nabla \left( \frac{1}{r} \right) = \frac{3}{r^3} \begin{bmatrix} x^2 & xy & xz \\ yx & y^2 & yz \\ zx & zy & z^2 \end{bmatrix} - \frac{1}{r^3} \begin{bmatrix} 1 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix} = \frac{3}{r^3} \begin{bmatrix} x^2 - \frac{1}{3} r^2 & xy & xz \\ yx & y^2 - \frac{1}{3} r^2 & yz \\ zx & zy & z^2 - \frac{1}{3} r^2 \end{bmatrix} \]

Now find the average gradient over the displacement:

\[ \left[ \nabla \left( \frac{1}{r} \right) \right]_{\text{avg}} \approx \frac{1}{2} \left[ \nabla \nabla \left( \frac{1}{r} \right) \right] = \frac{3}{2 r^3} \begin{bmatrix} x^2 - \frac{1}{3} r^2 & xy & xz \\ yx & y^2 - \frac{1}{3} r^2 & yz \\ zx & zy & z^2 - \frac{1}{3} r^2 \end{bmatrix} \]

\[ \Delta \Phi \approx Q \left[ \left( \frac{1}{r} \right) \right]_{\text{avg}} \cdot \mathbf{r} = Q \frac{3}{2 r^5} \left[ x^2 - \frac{1}{3} r^2 \right] \left( y^2 + xy' + xz' \right) 

+ yxx' + \left( x^2 - \frac{1}{3} r^2 \right) \left( y^2 + yz' + zxx' + zyy' + zzz' \right) + \left( z^2 - \frac{1}{3} r^2 \right) \left( z^2 - \frac{1}{3} r^2 \right) \]

There’s a trick here that’s kind of complicated that I don’t have time to write down yet, but it turns out we can drop the \((1/3)r^2\) terms in the first matrix by adding \(-\frac{1}{3} r^2\) terms to the quadrupole tensor. Briefly, it’s because \(\nabla \cdot \nabla \frac{1}{r}\) is traceless, so we can add any multiple of the identity matrix to the quadrupole tensor \(D_{\alpha\beta}\), and it won’t affect \(\Delta \Phi\). So we add \(-\frac{1}{3} \mathbf{I}\) to make \(D_{\alpha\beta}\) traceless. This means we can now add any multiple of the identity matrix to \(\nabla \cdot \nabla \frac{1}{r}\), and it won’t affect \(\Delta \Phi\). So we add \((r^2/3)\mathbf{I}\), to eliminate those ugly terms.

Now notice that all the terms for \(\Delta \Phi\) separate cleanly into \(r\) factors and \(r'\) factors. We also use the principle of superposition, to include the effect of all charges in the blob. This leaves us with
\[ \Delta \Phi_Q \approx Q_i \frac{3}{2r^5} \begin{bmatrix} x^2 & xy & xz \\ yx & y^2 & yz \\ zx & zy & z^2 \end{bmatrix} \text{element by element} \begin{bmatrix} \begin{bmatrix} x^2 - \frac{1}{3} r_i^2 & 0 & 0 \\ 0 & y^2 - \frac{1}{3} r_i^2 & 0 \\ 0 & 0 & z^2 - \frac{1}{3} r_i^2 \end{bmatrix} \end{bmatrix} \text{multiplied and} \begin{bmatrix} x' y' & x' z' \\ y' x' & y' z' \\ z' x' & z' y' \end{bmatrix} \text{summed with} \] 

So we define:

\[ D_{\alpha\beta} = \sum_{i=1}^{n} Q_i \begin{bmatrix} x_i' y_i' & x_i' z_i' \\ y_i' x_i' & y_i' z_i' \\ z_i' x_i' & z_i' y_i' \end{bmatrix} \begin{bmatrix} \begin{bmatrix} x_i^2 - \frac{1}{3} r_i^2 & 0 & 0 \\ 0 & y_i^2 - \frac{1}{3} r_i^2 & 0 \\ 0 & 0 & z_i^2 - \frac{1}{3} r_i^2 \end{bmatrix} \end{bmatrix} \]

and then:

\[ \Delta \Phi \approx \frac{3}{2r^5} \sum_{\alpha, \beta=1}^{3} D_{\alpha\beta} r_\alpha r_\beta = \frac{3}{2r^3} \sum_{\alpha, \beta=1}^{3} D_{\alpha\beta} \hat{r}_\alpha \hat{r}_\beta, \quad \text{where} \quad \hat{r}_\alpha = \alpha^{th} \text{ component of } \hat{r}. \]

\(D_{\alpha\beta}\) is the **quadrupole tensor**, which can simply be thought of as the matrix of numbers needed to compute the 2nd order change in \(\Phi\) due to a blob of charges. Notice that when summed with components of \(r\)-hat (the unit vector in the \(r\) direction), the change in \(\Phi\) varies as \(1/r^3\).

[Notice that the quadrupole tensor is a rank-2 tensor, which acts directly on another rank-2 tensor, to produce a scalar: \(\Delta \Phi\). The quadrupole tensor acts on the tensor \(r \otimes r\), which is the tensor product of the displacement vector with itself. We have thus seen two ways rank-2 tensors can act: some can act on a vector to produce a vector, and others can act on another rank-2 tensor to produce a scalar. It is not always sensible for a single rank-2 tensor to act both ways. See Funky Mathematical Physics Concepts for details.]

[Aside: The above last few steps would be a lot easier if we recognize that:

\[ (x, y, z) \begin{bmatrix} a & b & c \\ d & e & f \\ g & h & i \end{bmatrix} (x, y, z) = (x, y, z) \otimes (x, y, z) \]

\[ = \begin{bmatrix} x^2 & xy & xz \\ yx & y^2 & yz \\ zx & zy & z^2 \end{bmatrix} \text{element by element} \begin{bmatrix} a & b & c \\ d & e & f \\ g & h & i \end{bmatrix} \text{multiplied and} \begin{bmatrix} x^2 & xy & xz \\ yx & y^2 & yz \\ zx & zy & z^2 \end{bmatrix} \text{summed with} \begin{bmatrix} a & b & c \\ d & e & f \\ g & h & i \end{bmatrix}, \]

where we've defined a matrix dot-product, producing a scalar.]
4 Relativistic Electromagnetics

Relativity is second only to statistics in conceptual difficulty, because it is so contrary to common experience. However, when approached systematically, it can be understood, eventually. Here we not only present valid reasoning, but also describe the failure of common invalid reasoning. Topics:

1. Valid frame of reference
2. Time dilation and length contraction
3. Transformation of E and B fields
4. Doppler effect
5. Transformation of electromagnetic intensity

Construction of a Valid Frame of Reference

How to construct a valid reference frame:

(1) Fill space with stationary clocks. Two clocks can be synchronized by putting a flashing light source halfway between them. Synchronize all your clocks; they remain synchronized.

(2) Fill space with a coordinate system. Measure the distances to vertices in the coordinate system with stationary measuring rods.

(3) Measure any event as its position on the coordinate system, and its time measured by a clock right at the event. Therefore, there is no propagation delay between the event and the clock, and no other funny relativistic effect, either.

Example: Measuring speed: The wrong way: we’re used to marking off miles on the road, driving over the road in our car, and timing the distance between mile markers. But this overstates our speed. Why? Because once we start moving in the car, the miles on the road shrink from length contraction, and we think we’re traveling farther than we really are. Remember, “really” means “distances as measured in our frame of reference,” i.e. measured by rods at rest with respect to us. No other method of measuring distance is valid. The right way: we measure our speed relative to the road by measuring how long it takes a single point on the road to pass between two known points in our reference frame. The two known points are stationary relative to us, and we use two clocks, one at each point, to measure the time interval. Since all our clocks are synchronized, we can always use different clocks to measure time at different events.

By symmetry, if we measure our speed relative to the road, and someone on the road measures our speed relative to the road, we must get the same answer. Since either frame is valid, there is no basis for one frame to measure faster than the other, so the two observers must measure the same speed.

TBS: distance perpendicular to the motion must be the same for both observers. [Helliwell p??].

Time Dilation and Length Contraction

Alice and Bob are moving relative to each other, but of course, each has a measuring system (reference frame) that is stationary relative to himself.
Bob travels in the +x direction (parallel to the wall) at speed $\beta = v/c$. What does Bob measures? First, Bob sees himself as stationary, and Alice as moving to the left. He sees that the light beam is not perpendicular to the wall, so we conclude that the angle at which a light beam travels is relative; i.e., different observers measure different angles. (This is true of baseballs or anything else that moves.) Also, Bob sees the light travel a larger distance than Alice does; since the speed of light is the same for Alice and Bob, Bob clocks a longer time between the 2 events. Call Bob’s time $\gamma$ (in seconds). During that time, Alice moves left a distance (by his measure) of $\beta \gamma$ light seconds. Since distances perpendicular to the motion are the same for Alice and Bob, Bob measures a $y$ displacement of 1 (light-second). Therefore, Bob’s time is:

$$\gamma^2 = (\beta \gamma)^2 + 1^2, \quad \gamma^2 - \beta^2 \gamma^2 = 1, \quad \gamma = \frac{1}{\sqrt{1 - \beta^2}}. $$

During this time, a clock stationary in Alice’s frame measures only 1 second. Therefore, Bob sees a moving clock as ticking more slowly than his own (stationary) clocks. This is \textbf{time dilation}. By symmetry, Alice (and everyone else in the universe) must see the same phenomenon.

Moving clocks tick more slowly than stationary ones.

This is \textit{not} an illusion caused by the finite speed of light (see discussion of valid reference frames above). This is real.

Do not be misled by this erroneous argument: Alice sees one second on her clock, and sees Bob time $\gamma > 1$ seconds on his clocks, therefore Alice thinks Bob’s moving clocks run faster. The error is that Bob used two clocks, at different $x$ positions, to make his measurement. In Bob’s frame, those clocks are synchronized, but in Alice’s frame they are not.

If Alice looks at any \textit{one} of Bob’s clocks (not two clocks in two different $x$ positions), then she will indeed see Bob’s clock run more slowly.

\textbf{Length contraction:} By symmetry, Alice and Bob agree that they are moving at a speed $\beta$ relative to each other. When Alice turns on her flashlight, she simultaneously marks a green dot on Bob’s moving frame. When she detects the light on the wall, she marks a red dot on Bob’s frame. During the interval, Alice measures Bob move $\beta \Delta t = \beta$ light-seconds across her frame; i.e. Alice measures the $x$ distance between two moving dots as $\beta$ light-seconds. The dots are stationary in Bob’s frame, and as already noted, Bob measures their separation as $\beta \gamma$ light-seconds. Therefore, Alice measures the length of moving things as shorter than someone at rest with respect to those things, by the factor $1/\gamma$.

Do not be misled by this erroneous argument: Look at the diagram of Alice’s measurement of the distance between the dots she painted on Bob’s reference frame. As noted, Alice measures that distance as $\beta$. Suppose Alice has a ruler on her ground, with which she measures that distance. One might (incorrectly) conclude that at that moment, Bob sees her moving ruler aligned with his dots, and conclude that in his frame, the ruler measures $\gamma \beta$, and thus moving rulers \textit{grow} instead of
contract. This is wrong because in Bob’s frame, the dots do not align with Alice’s ruler at a single instant in time.

For the correct reasoning, see Figure 4.1d: In Bob’s frame, Alice is moving to the left, and her ruler is short. Bob measures the red dot aligns with the red end of Alice’s ruler after the green dot aligns with the green end of Alice’s ruler.

Moving objects are measured shorter than when at rest.

Transformation of E & B Fields

There’s a simple picture of the relativistic transformation of E & B fields. We seek \( E' \) and \( B' \), the fields in the moving frame, in terms of \( E \) and \( B \), the fields in the stationary (lab) frame. First, we’ll find \( E' \), the E-field in the moving frame. We start with transformation of the E-field alone. Imagine an E-field created by a sheet charge:

\[
\sigma' = \gamma \sigma \Rightarrow E_\perp' = \gamma E_\perp
\]

In the moving frame, the sheet charge is compressed to a smaller area, by the length contraction factor \( \gamma \). Therefore, \( E' \) is increased by the same factor.

Now we add a B-field to the stationary frame:

\[
E_\perp' = \gamma (E_\perp + v \times B_\perp) \quad (\text{SI})
\]

A stationary charge in the moving frame, is (in the lab frame) moving relative to \( B \). Nonrelativistically, the charge feels a force of \( F = qv \times B \), just like an E-field of \( v \times B \) (SI units), simply from the Lorentz force law. This appears in the moving frame to be an E-field: \( E' = v \times B \). Boosting to relativistic speeds, picture that the E-field can be thought of as “lines of force” whose density per unit area is proportional to the field strength. Because of Lorentz contraction, the same lines of force occupy a smaller area in the moving frame, and so their density is increased by \( \gamma \):

\[
E_\perp' = \gamma (E_\perp + v \times B_\perp), \quad E_\parallel' = E_\parallel \quad [\text{Jac 11.148 p558}].
\]

There is no contraction of the plane perpendicular to the direction of motion, and no Lorentz force from \( B_\parallel \) (parallel to the velocity). Hence, \( E_\parallel' = E_\parallel \).

\( B' \) is similar. TBS??

\[
B_\perp' = \gamma (B_\perp - v \times E_\perp), \quad B_\parallel' = B_\parallel \quad [\text{Jac 11.148 p558}].
\]
We can combine the $E$ (and $B$) transformations above into a single vector equation for each. If we start with the perpendicular formulas, and note that the cross-product term does not contribute anything to the parallel component, then replacing $B_\perp \rightarrow B$ would incorrectly inflate the parallel component by $\gamma$. We can fix this by just subtracting off $(\gamma-1)\hat{\beta} \cdot E$:

$$E' = \gamma(E + \beta \times B) - (\gamma-1)\hat{\beta} \cdot E \quad \quad \quad B' = \gamma(B - \beta \times E) - (\gamma-1)\hat{\beta} \cdot B.$$ 

Jackson complicates this formula with the identity $(\gamma-1)\hat{\beta} \cdot E = \frac{\gamma^2}{\gamma+1} \beta (\beta \cdot E) \Rightarrow$

$$E' = \gamma(E + \beta \times B) - \frac{\gamma^2}{\gamma+1} \beta (\beta \cdot E), \quad B' = \gamma(B - \beta \times E) - \frac{\gamma^2}{\gamma+1} \beta (\beta \cdot B) \quad \text{[Jac 11.149 p558]}.$$

### Acceleration Without Force

Consider a particle moving in the $x$-direction, and a force pushing it in the $+y$-direction. The particle accelerates up, but it also decelerates in the $x$-direction, even though there is no force in the $x$-direction. That’s because the acceleration in $y$ increases the particle’s magnitude of velocity (speed), and therefore the particle’s $\gamma = \frac{1}{\sqrt{1 - v^2/c^2}}$. The $x$-momentum doesn’t change, but when $\gamma$ increases, $v_x$ must decrease to keep the same momentum:

$$F_y = \frac{dp_y}{dt} \quad \Rightarrow \quad v_y \text{ increases, and } |v| \text{ and } v^2 \text{ increase.}$$

$$F_y = 0 = \frac{dp_y}{dt} \quad \Rightarrow \quad p_y = \gamma mv_y = \text{const.} \quad \gamma \text{ increasing} \quad \Rightarrow \quad v_y \text{ decreases!}$$

### On-Axis Doppler

We first consider the simpler case of motion along the line of propagation. We are a stationary light source, and there is a moving observer traveling directly toward us (above). We can deduce what the observer sees from our own observations. Note that wave crests are countable things, and therefore the same in all frames. In an interval of our time $dt$, the observer starts at A and ends at B, and simultaneously, the wave train starts at C and also ends at B. During this interval, the observer sees all the crests between A and C, i.e.

$$n = \left( f \frac{dt}{c} \right) \left( \frac{c + v}{c} \right) = \left( f \frac{dt}{c} \right) \left( 1 + \frac{v}{c} \right) \quad \text{where} \quad v \text{ is the speed of approach}.$$ 

However, his clock advances only $(dt / \gamma)$ seconds, due to time dilation. Without loss of generality, let $dt = 1 \text{ second}$. Then in one of the observer’s seconds, he sees

$$n' = f' = \gamma f \left( 1 + \frac{v}{c} \right).$$

This formula is valid for positive and negative $v$. And of course, only the relative motion of the source and observer matter (after all, this is relativity, isn’t it?)
Off-Axis Doppler

For variety, we now take the reverse (but equivalent) view that you are now the observer, you are stationary, and the light source is moving. (I think this is a little harder to follow). You are in the enterprise, and a star has an off-axis relative motion of \( v \) (see diagram).

Let \( r = \text{distance from enterprise to star} \)
\( x = \text{“horizontal” distance to star} \)
\( v = \text{speed of star in x direction} \quad (v < 0) \)
\( f_s = \text{frequency at which star emits light (in its own frame)} \)
\( f_o = \text{frequency at which enterprise sees light} \)

From the diagram, and recalling that \( v < 0 \),
\[
\frac{dr}{dt} = 2x \frac{dx}{dt} = 2r \cos \theta v
\]
\[
dr = v (\cos \theta) \ dt
\]

Because of time dilation, the emitted frequency in the Enterprise frame is multiplied by a factor of \( (1/\gamma) \). In addition, the waves are “compressed” by the changing separation between the star and Enterprise. In a time interval \( dt \), the Enterprise sees the star emit \( (f_s/\gamma) \ dt \) cycles of EM wave. But those cycles get squeezed into a space of \( (c \ dt + dr) \) distance \( (dr < 0) \). Thus, the number of cycles the enterprise sees in a full distance of \( (c \ dt) \) is (recalling again that \( v < 0 \)):
\[
f_o \ dt = \frac{f_s}{\gamma} \left( \frac{c \ dt}{c \ dt + dr} \right) = \frac{f_s}{\gamma} \left( \frac{c \ dt}{c \ dt + v \cos \theta \ dt} \right)
\]
\[
f_o = \frac{f_s}{\gamma} \left( \frac{c}{c + v \cos \theta} \right) = \frac{f_s}{\gamma (1 + \beta \cos \theta)}
\]

This reduces to the well known simpler formula for on-axis motion when \( \theta = 0 \):
\[
f_o = \frac{f_s}{\gamma (1 + \beta \cos \theta)} = \frac{f_s}{\gamma (1 + \beta)}.
\]

Transformation of Intensity (Power Density)

If a light wave has an intensity \( I \) in the stationary frame, what is its intensity \( I' \) as seen in an on-axis moving frame? We first find this from the transformation of the E field, and noting that \( I \propto E^2 \). Then we show another, more intuitive, way to find the intensity transformation.

First, the transforming E-field approach: In gaussian units, a free space wave as \( B = E \). For on-axis motion:
\[
E' = \gamma \left( E + \frac{v}{c} \times B \right) \quad \Rightarrow \quad E' = \gamma \left( E + \frac{v}{c} B \right) = \gamma \left( E + \frac{v}{c} E \right) = E \gamma \left( 1 + \frac{v}{c} \right).
\]

Electric fields in propagating waves transform exactly like the Doppler frequency shift.

Then, for source and observer approaching:
\[
I \propto E^2 \quad \Rightarrow \quad I' = \gamma^2 \left( 1 + \frac{v}{c} \right)^2 I = \frac{(1 + \frac{v}{c})^2}{(1 - \frac{v}{c})(1 + \frac{v}{c})} = \frac{1 + \frac{v}{c}}{1 - \frac{v}{c}} I.
\]
Intensity transforms as the square of the Doppler shift.

It is perhaps more intuitive (though less theoretically sound) to consider the light wave as a “flux of photons,” i.e. some number of photons per unit area per unit time, impinging on the observer. The flux times the photon energy is proportional to intensity. If the flux seen by a stationary observer is $f$, what is the flux seen by a moving observer, $f'$? The picture is essentially identical to that we used for deriving on-axis Doppler shift:

Without loss of generality (wlog), the tube has cross-sectional area of $1\text{m}^2$. As seen by you (stationary), the moving observer intercepts $n_t = \left(f \frac{dt}{\gamma} \right) \left(1 + \frac{v}{c}\right)$ photons in the time interval $dt$. But again, his clock only advanced $(dt/\gamma)$ seconds. Therefore, the observer’s photons per unit time (the cross sectional area for him is still $1\text{m}^2$) is $n' = f' = \gamma f \left(1 + \frac{v}{c}\right)$. The flux transforms like Doppler shift. (Note also that as the moving observer looks ahead at the photon “density” in space, his density is greater than yours by a factor of $\gamma$, because the tube is shorter by a factor of $\gamma$ due to length contraction.)

But each photon’s energy is proportional to its frequency, $E = \hbar \omega$, and each photon’s frequency is Doppler shifted by $\gamma (1 + v/c)$, as shown earlier. Ergo,

$$I' \propto n' E' \quad \Rightarrow \quad I' = I \gamma \left(1 + \frac{v}{c}\right) \gamma \left(1 + \frac{v}{c}\right) = I \gamma^2 \left(1 + \frac{v}{c}\right)^2 \quad v \equiv \text{approach speed},$$

as before.

**How Big Is a Photon?**

So what’s the theoretical objection to the above photon-flux model of intensity transformation? The problem is that photons are not point-like particles of small dimensions. The electromagnetic wave is a widely distributed wave-function (of sorts) of the photon, in the quantum mechanical sense. However, in quantum field theory (the only valid quantum theory for photons), the photon density is proportional to the square of the electromagnetic intensity, just as with massive particles. The photon itself, though, cannot be localized, or thought of as a point, but the density approach is rigorously valid.

There are many real-world situations in which a single photon is spread over many meters or kilometers, e.g., lunar laser ranging which detects individual photons spread laterally over many square meters of the receiving telescope.

**The Pressure of Light**

We can derive the pressure of light from the fundamental relativistic energy momentum relationship:

$$E^2 = \left(mc^2\right)^2 + \left(pc\right)^2$$

Since light is (i.e., photons are) massless, we have:

$$E = pc \quad \text{(light or other massless particles)}$$
Now force  \( F = dp/dt \), so for normal incidence light absorbed onto a surface:

\[
F = \frac{dp}{dt} = \frac{1}{c} \frac{dE}{dt} = \frac{P}{c}
\]

where \( P \equiv \text{power of radiation} \)

Taken per unit area:

\[
\text{Pressure} = \frac{I}{c}
\]

where \( I \equiv \text{intensity} = \text{power per unit area} \)

**Example: Reflection Off a Moving Mirror**

We show here a novel solution to the classic test question of light reflecting off a moving mirror, using conservation of energy entirely in the lab frame.

**The question:** An ideal mirror moves to the right, perpendicularly to its face. A physicist shines a light on it with normal incidence from the left (Figure 4.2). Find the power reflection coefficient of the moving mirror \( \Gamma \equiv (P_r/P_i) \).

**The intensity transformation way:** In previous sections, we derived (in two different ways) that \( E \) and \( B \) fields transform like the Doppler shift, and therefore intensity transforms like the square of Doppler (the minus sign below is in the numerator because source and observer are receding):

\[
I' = I_i \left( \frac{1-v/c}{1+v/c} \right) \quad v \equiv \text{recession speed}.
\]

In this problem, we have two intensity transformations of the same factor: once into the mirror frame, and once again from the mirror frame to the lab frame. The result is then simply

\[
I_r = I_i \left( \frac{1-v/c}{1+v/c} \right) = I_i \left( \frac{1-v/c}{1+v/c} \right)^2 \quad \Rightarrow \quad \Gamma = \left( \frac{1-v/c}{1+v/c} \right)^2.
\]

**The energy way:** It is perhaps instructive to consider conservation of energy in the lab frame: the energy crossing a transverse plane (parallel to the mirror) goes into (1) filling up the increasing distance to the mirror with electromagnetic energy, (2) doing work on the mirror from the pressure of light, and (3) transmitting light across the boundary back into the laboratory. Without loss of generality, we can use an incident power of 1 (in arbitrary units). In some time interval, the fraction of the incident energy that gets stored in the expanding space up to the mirror is \( v/c \). The rest of the light hits the mirror, and reflects. Similarly, the fraction of reflected energy that gets stored in the expanding space is \( \Gamma v/c \).

The work done on the mirror by the incident light is \( F \Delta x \), where \( F = P/c \), and \( \Delta x \) in a unit time is \( v \). Then \( W = P v/c \). However, \( P \) is not \( P_i = 1 \), but just the fraction of \( P_i \) that is not stored in the expanding volume. We found above that the stored energy is \( v/c \), so that which hits the mirror is \( (1-v/c) \), and the work is \( W = (1-v/c) v/c \).

Similarly for the work of the reflected light, but the total work is from the light stored in the expanding volume plus the light sent back through the transverse plane into the lab. The stored energy is \( \Gamma v/c \), and the reflected light power is \( \Gamma \). Hence the work is \( W = \Gamma (1+v/c) v/c \).

Then by conservation of energy, the incident energy equals the sum of stored incident plus reflected energy, plus the incident plus reflected work:

\[
W_{\text{incident}} = W_{\text{incident stored}} + W_{\text{reflected}} + W_{\text{incident work}} + W_{\text{reflected work}}.
\]
\[ 1 = \frac{v}{c} + \left( \frac{1-v}{c} \right) \left( \frac{v}{c} \right) + \Gamma \left( \frac{1+v}{c} \right) \left( \frac{v}{c} \right) + \Gamma \] 

\[ \Rightarrow 1 - 2 \frac{v}{c} = \Gamma \left( 1 + 2 \frac{v}{c} + \left( \frac{v}{c} \right)^2 \right) \]

\[ \Gamma = \left( \frac{1-v/c}{1+v/c} \right)^2 \]

**Beaming**

Both Jackson and Griffiths have misleading diagrams for relativistic beaming at high \( \gamma \). Both their diagrams show no backward emission components at all. Inspection of the equation shows immediately that this is not the case; the power is finite for all angles up to \( \pi \). Here are more suggestive diagrams:

![Diagram](image)

**Figure 4.3** Radiation pattern (cross-section) for transverse acceleration at various velocities.

Remember that the radiation pattern is axially symmetric about the velocity vector, so the “beaming” is more like a big hollow cone of light than a pair of headlights.

**Covariant Form of Maxwell’s Equations**

Given that \( F^{\mu\nu} \) is a rank-2 tensor, we can show that \( J^\mu \) is a 4-vector:

Define \( \Lambda^\mu_\delta = \frac{\partial x^\mu}{\partial \tilde{x}^\delta} \) = Lorentz transformation matrix from \( s \) to \( \tilde{x} \)

\[ \lambda_\nu^\alpha = \frac{\partial \tilde{x}^\alpha}{\partial x^\nu} = \left( \Lambda^\mu_\delta \right)^{-1} \] = Lorentz transformation matrix from \( \tilde{x} \) to \( s \)

Then

\[ \lambda_\nu^\alpha \lambda_\nu^\mu = \delta^\alpha_\mu \] the identity transformation

Note \( \partial_\mu = \lambda^\nu_\mu \partial^\nu \) or from \( \frac{\partial}{\partial \tilde{x}^\nu} = \frac{\partial \tilde{x}^\nu}{\partial x^\mu} \frac{\partial}{\partial x^\mu} \)

In \( s \):

\[ \partial_\nu F^{\mu\nu} = \mu_0 J^\mu \]

In \( \tilde{x} \):

\[ \partial_\nu \tilde{F}^{\mu\nu} = \frac{\partial \tilde{F}^{\mu\nu}}{\partial \tilde{x}^\nu} = \partial_\alpha \lambda^\mu_\alpha \lambda^\nu_\delta F^{\delta\lambda} = \Lambda^\mu_\delta \partial_\alpha F^{\delta\alpha} = \mu_0 \Lambda^\mu_\delta J^\delta = \mu_0 J^\mu \]
5 Shorts

Is the Electric Field Real?

Is the electric field real? Or is it just a tool to help model reality? This question has been asked over the history of our understanding of EM fields, and the answer has evolved over time. For the electric field:

First, people noticed that certain particles pushed or pulled on each other. To model that, they invented the idea of “charge”, and Coulomb’s force law. At this point, there is no need for any fields, since electric forces are all two-particle interactions.

In more complicated situations, particularly when there is a fixed set of charges, called the “source” charges, it was convenient to pre-compute the effect of these source charges on hypothetical “test” charges. This pre-computation worked because of the linearity of the Coulomb force. The result was a vector function of space, and was given the name “electric field.” At this point, the electric field is purely a mathematical convenience for describing point-charge interactions, with no physical significance. [Note that all forces, by the definition of a vector, must add.]

Later, experiments showed that if energy is to be conserved, then it is convenient to say that this “electric field” actually stores energy. Now the electric field starts to seem like a physical reality.

Still later, experiments showed that a time varying electric field could produce waves which propagate very far away from any source charges. Essentially, the electric field (together with the magnetic field) is self-sustaining. This makes fields seem quite real indeed, as they are now far separated from their sources, and seem to exist independently of them. This view is somewhat unsatisfactory, though, because the ultimate source of all EM radiation is still accelerating charges.

However, this time-dependent theory of EM fields also leads to the experimental fact that the effect of a charge on a distant charge is delayed by a propagation time. What is propagating? If fields are real entities, then the answer is clear.

Furthermore, when a particle radiates, it recoils immediately, in a direction opposite to the radiation. This implies a transfer of momentum from the particle to the radiation, which is immediate, and independent of when that radiation might later be absorbed. Thus, the energy and momentum of EM radiation appears to be quite “real,” and the EM field’s energy and momentum exist independently of the particles.

The final evidence favoring fields as “real” is that they contribute to the center of mass, now called center of energy, of a system of particles. This goes beyond conservation of energy, which counts only the total energy, because now the location of the energy in space matters. To satisfy the condition that the center of energy remain fixed for an isolated system, we must associate an energy density with the electric and magnetic fields, such that:

\[ \rho_{\text{electric}} \propto E^2, \quad \rho_{\text{magnetic}} \propto B^2. \]

The currently accepted classical (non-quantum) model is that fields are real things: they carry both energy and momentum, and we can say where the energy and momentum is located, e.g., the energy density varies as the square of the field.

Quantum mechanics is fully consistent with this classical picture, as photons (quantized EM fields) are treated fully and equally with matter as particles: they carry both energy and momentum.

The complete picture is somewhat more complicated due to the details of magnetic fields, and their interaction with electric fields. However, the above reasoning and conclusions remain valid.

The Coulomb Force Constant is Defined Exactly

Recall that \( c \) and \( \mu_0 \) are defined exactly:
Charge Is the Fundamental Electrical Quantity

Conceptually, there are 4 fundamental macroscopic quantities in the universe: distance, time, mass, and charge. This is a conceptual truth that does not depend on how we choose our units of measure. For example, in the MKSA (aka SI) system, the unit of charge is the coulomb. However, as a matter of laboratory convenience, the SI system defines the ampere first, and then defines the coulomb as the amount of charge flowing past a point in 1 second, with a current of 1 ampere. This is simply because it’s easier to measure magnetic forces in the lab than electrostatic forces. This does not change the theoretical reality that charge is one of the 4 fundamental measurable physical quantities. The ampere is a compound quantity involving charge and time. Particles have charge; they don’t have currents.

In contrast to SI, note that CGS units define the charge as the first quantity, and current in terms of charge. Clearly, this variety of human chosen definitions has no bearing on the nature of physics.

Units of Measure: SI vs. Gaussian

When converting equations between SI and Gaussian units, three simple observations will help tremendously in remembering the conversions. Three formulas are easy to remember in both units, since they are so common: (1) electric potential, (2) the speed of light, and (3) the Lorentz magnetic force. These provide all the information needed for most conversions. By comparing these equations in each set of units, we can see what substitutions are needed to take almost any equation from one system to the other. We show here conversion from SI to Gaussian, because undergraduates usually learn in SI, then have to switch to Gaussian as graduates:

<table>
<thead>
<tr>
<th>Notes</th>
<th>SI</th>
<th>Gauss</th>
</tr>
</thead>
<tbody>
<tr>
<td>(1)</td>
<td>( \phi(r) = q / 4\pi\varepsilon_0 r )</td>
<td>( \phi(r) = q / r )</td>
</tr>
<tr>
<td>(2)</td>
<td>( c^2\varepsilon_0\mu_0 = 1 )</td>
<td>( \varepsilon_0\mu_0 = 1 )</td>
</tr>
<tr>
<td>(3)</td>
<td>( \mathbf{F} = q\mathbf{v} \times \mathbf{B} )</td>
<td>( \mathbf{F} = q \frac{\mathbf{v}}{c} \times \mathbf{B} )</td>
</tr>
</tbody>
</table>

(1) This does not mean that \( \varepsilon_0 \) is \( 1/4\pi \) in Gaussian units; in fact, \( \varepsilon_0 = 1 \) in Gaussian units. It means that when including the whole package of conversions to take an equation from SI to Gaussian, including the unit of charge, the final result is that \( \varepsilon_0 \) turns into \( 1/4\pi \).

(2) Note that we used #1 above \( (\varepsilon_0 \rightarrow 1/4\pi) \) to eliminate \( \varepsilon_0 \).

(3) This means that the B-field in Gaussian units is inflated by the factor \( c \), so when using it, you have to deflate it to make it work in the equation. For example, that’s why a propagating wave has:

\[
\text{(SI)} \quad E = cB, \quad \text{but} \quad \text{(gaussian)} \quad E = B.
\]

Example: Biot-Savart Law:

\[
\text{(SI)} \quad dB = \frac{\mu_0}{4\pi} \frac{Idl \times \hat{r}}{r^2} \rightarrow \quad \text{(gaussian)} \quad \frac{dB}{c} = \frac{4\pi}{4\pi c^2} \frac{Idl \times \hat{r}}{r^2} \Rightarrow \quad dB = \frac{1}{c} \frac{Idl \times \hat{r}}{r^2}.
\]

Bound and Gagged

Are bound charges real? To understand this, we must be more specific: there are real bound charges that are responsible for permittivity (and equivalently, for dielectric constant). However:
The so-called “bound charge distributions” that one can calculate are not real (despite what some references say).

A simple example proves the point. Consider a broad (i.e., infinite) flat sheet of dielectric material:

![Diagram showing dielectric with and without an electric field](image)

**Figure 5.1** (Left) Dielectric with no E-field. (Middle) Dielectric with long dipoles in an E-field. (Right) Dielectric with higher volume density of shorter dipoles in E-field.

The diagram shows only a representative volume, but the top and bottom sheets are given as broad. Inside the dielectric are molecules with electric dipole moments. Above left, absent an E-field, the dipoles are randomly oriented. They therefore sum to zero total dipole moment. Above middle, in an E-field, the dipoles align. In this example, the dipoles are long, and the positive (blue) ends of the top two rows land on top of, and cancel, the negative (red) ends of the bottom two rows. The net result is that all inside charges cancel, and we are left with two surface charges, negative on top, and positive on the bottom. These charge densities are what one computes with the standard formula

\[ \sigma_{\text{bound}} = \mathbf{P} \cdot \mathbf{n} \]

The electric field from the dipoles is uniform throughout the slab.

However, this is clearly an unrealistic depiction. Real molecules aren’t that close together; their dipole moments are not ideal point charges separated by vacuum; and adjacent charges don’t exactly cancel.

More realistically (but still idealized), above right, the dipoles don’t overlap (they’re shorter); each one has a larger dipole moment, so the endpoint charges are larger; and there are more of them. However, the dipole moment per unit volume is the same as the middle diagram. Therefore, these dipoles add up to the same total dipole moment in the slab as the middle diagram, but they result in larger real surface charge densities. In addition, there are sheet charges at other layers below the surface, which cause the induced E-field to be large between the charges of each dipole, and zero in the gaps between layers of dipole. That’s because a real dipole moment is spread throughout the volume of the dielectric. In fact, the standard definition of polarization for a dielectric (in an E-field) is dipole moment density (dipole moment per unit volume). The total dipole moment in the slab is the product of moment-density times volume:

\[ \mathbf{p} = (\mathbf{P})(\text{Volume}) \]

where \( \mathbf{p} = \text{dipole moment}; \ \mathbf{P} = \text{dipole moment density} \).

Therefore:

The bound surface charge densities we compute for dielectrics are model densities which give equivalent dipole moment to the real dielectric, but are not real.

**Pole to Pole: Current Loops Are Approximate Dipole Moments**

A current loop has mostly a dipole moment, but any finite-sized loop also has other multipole moments in it. A pure dipole results only in the limit that the loop size goes to zero, while the current increases to infinity, in such a way that the product \( IA \) remains constant. (This is similar to an electric dipole, which is a pure dipole only in the limit of the charge \( q \to \infty \), while separation \( d \to 0 \), such that the product \( qd = \) constant.) For example, in a pure magnetic dipole pointing up, all the magnetic field in the equatorial plane points down:

\[ \mathbf{B}(\mathbf{r}) = \frac{3(\mathbf{M} \cdot \mathbf{\hat{r}}) \mathbf{\hat{r}} - \mathbf{M}}{r^3} \]

[Jac 5.56 p186, Gri 3.103 p153]
But in a finite loop, the field inside the loop points up. Hence, a finite loop cannot be a pure dipole.

\[ F \equiv \frac{d\mathbf{p}}{dt} = q \left( \mathbf{E} + \frac{\mathbf{v}}{c} \times \mathbf{B} \right) \]

\( q \equiv \text{test charge. gaussian units, relativistically valid} \).

This says that the \( \mathbf{E} \) field acts on any charge, moving or not. The \( \mathbf{B} \) field acts only on moving charges. The \( \mathbf{E} \) field can do work on a charge, but the \( \mathbf{B} \) field cannot directly do work on a charge. The \( \mathbf{E} \) and \( \mathbf{B} \) fields are directly measurable by their effects on a test charge. Physically, the sources of \( \mathbf{E} \) fields are (1) charges (moving or not), and (2) changing \( \mathbf{B} \) fields. The source of \( \mathbf{B} \) fields is charge currents (i.e., moving charges).

Sometimes however, instead of working directly with \( \mathbf{E} \) and \( \mathbf{B} \) fields, it is more convenient to work with potentials, which generate the \( \mathbf{E} \) and \( \mathbf{B} \) fields by a simple mathematical operation. Two such potentials are needed: a scalar potential \( \Phi(t, \mathbf{x}) \), and a vector potential \( \mathbf{A}(t, \mathbf{x}) \). These potential functions are not unique: different choices can produce the exact same \( \mathbf{E} \) and \( \mathbf{B} \) fields, and thus the same physical predictions.

Static Fields

The defining property of the magnetic vector potential \( \mathbf{A}(t, \mathbf{x}) \) is that \( \mathbf{B}(t, \mathbf{x}) = \nabla \times \mathbf{A}(t, \mathbf{x}) \). Static fields are not functions of time, so in the static case, the above definition can be reduced to:
\[ \mathbf{B}(\mathbf{x}) = \nabla \times \mathbf{A}(\mathbf{x}) \]  

Then we can define an electric potential \( V(\mathbf{x}) \) that directly relates to energy:

\[
\text{Define } V(\mathbf{x}) \text{ such that } \mathbf{E}(\mathbf{x}) = -\nabla V(\mathbf{x}) \implies V(\mathbf{x}) = -\int_{x_0}^{x} \mathbf{E}(\mathbf{x}) \cdot d\mathbf{x}.
\]

where we choose \( V(\mathbf{x}_0) = 0 \).

This electric potential has no gauge freedom (only an arbitrary additive constant). The \( \mathbf{E} \)-field is given by the gradient of the electric potential, which in the static case also serves as a scalar potential:

\[
\mathbf{E}(\mathbf{x}) = -\nabla V(\mathbf{x}) = -\nabla \Phi(\mathbf{x}) \quad \text{where} \quad V(\mathbf{x}) \equiv \text{electric potential} \quad \text{(static fields)}.
\]

\( V(\mathbf{x}) \) has units of joules/coulomb, or volts. Both of these equations involve only derivatives of the potentials, so adding a constant to \( \Phi \) or \( \mathbf{A} \) doesn't change the fields. This is a trivial form of gauge invariance, which we ignore.

More importantly, the \( \mathbf{A} \)-field is not unique because there are many functions \( \mathbf{A}(\mathbf{x}) \) which have the same curl. In particular, any irrotational (i.e., curl-free) vector field can be added to \( \mathbf{A} \), and it doesn't change the curl.

Any irrotational field can be written as the gradient of a scalar field, and the gradient of any scalar field is irrotational.

Our gauge freedom for static fields then means we can add the gradient of any scalar function to \( \mathbf{A} \), and get another, physically equivalent, vector potential:

\[
\mathbf{A}'(\mathbf{x}) = \mathbf{A}(\mathbf{x}) + \nabla \Lambda(\mathbf{x}),
\]

where \( \mathbf{A}' \) is physically equivalent to \( \mathbf{A} \), and \( \Lambda(\mathbf{x}) \) is any scalar function.

**Time Varying (Dynamic) Fields**

Things get more complicated in the dynamic case. Following [Jac sec. 6.2-6.5 p239-48], electric and magnetic fields can still be specified in terms of a scalar potential and vector potential, but now a time-varying \( \mathbf{B} \)-field, and thus a time-varying \( \mathbf{A} \)-field, contributes to the \( \mathbf{E} \)-field:

\[
\nabla \times \mathbf{E}(t, \mathbf{x}) = -\frac{\partial \mathbf{B}(t, \mathbf{x})}{\partial t} = -\frac{\partial}{\partial t} \left( \nabla \times \mathbf{A}(t, \mathbf{x}) \right) \quad \Rightarrow \quad \nabla \times \left( \mathbf{E}(t, \mathbf{x}) + \frac{\partial \mathbf{A}(t, \mathbf{x})}{\partial t} \right) = 0.
\]

Note that the curl and time derivatives commute, so we can bring the \( \partial / \partial t \) inside the curl operator. Since the curl of \( \mathbf{E} + \partial \mathbf{A}/\partial t \) is zero, it can be written as the gradient of a scalar function which we call \( -\Phi \):

\[
\mathbf{E}(t, \mathbf{x}) + \frac{\partial \mathbf{A}(t, \mathbf{x})}{\partial t} = -\nabla \Phi(t, \mathbf{x}) \quad \text{and} \quad \mathbf{B}(t, \mathbf{x}) = \nabla \times \mathbf{A}(t, \mathbf{x}).
\]

For dynamic fields, \( \Phi(t, \mathbf{x}) \) is not the electric potential. It is the scalar potential, chosen for mathematical convenience, which, along with a matching \( \mathbf{A}(t, \mathbf{x}) \), generates \( \mathbf{E}(t, \mathbf{x}) \).

In fact, the electric potential does not exist (in general) in the time varying case, since the electric field has a non-zero curl. This means that the energy \( \Delta E \) required to move a charge from a point \( \mathbf{x}_1 \) to \( \mathbf{x}_2 \) depends on the path, and hence there cannot exist any function \( V(\mathbf{x}) \) such that \( \Delta E = V(\mathbf{x}_2) - V(\mathbf{x}_1) \). Nonetheless, it is possible to define the electrostatic potential, which accounts for only the coulomb forces of source charges, but not the dynamics of varying \( \mathbf{B} \)-fields.

Since both \( \mathbf{E} \) and \( \mathbf{B} \) are found from derivatives of \( \mathbf{A} \) and \( \Phi \), \( \mathbf{A} \) and \( \Phi \) are not unique. **Gauge freedom** (or **gauge invariance**) is the ability to impose additional mathematical constraints on \( \mathbf{A} \) and \( \Phi \), that don't
change the physical system. [It is sometimes said that $E$ and $B$ are the “physical” fields, and $A$ is not; this point is debatable, but the preceding definition of gauge invariance holds either way.]

**Lorenz Family of Gauges**

You may hear talk of “the” Lorenz gauge (often incorrectly referred to as “Lorentz” gauge [Jac p294]). Lorenz gauges are actually a family of gauges, which satisfy

$$\nabla \cdot A(t, x) + \frac{1}{c^2} \frac{\partial \Phi(t, x)}{\partial t} = 0 \quad \text{Lorenz gauge} \quad \text{[Jac 6.14 p240]}$$

or

$$\partial_\mu A^\mu(t, x) = 0 \quad \text{Relativistic form} \quad \text{[Jac 11.133 p555]}$$

[Aside: As a simple example of gauge invariance, look at the first equation above. We can generate another Lorenz gauge vector potential $A'(t, x)$ by adding any static (i.e., time-independent) divergenceless vector field $G(x)$ to a given Lorenz gauge field $A$. Since the curl of $G$ must be 0 (to preserve the $B$-field), $G$ can be written as the gradient of a scalar function: $G(t, x) = \nabla \Lambda(t, x)$.]

Since we chose $\nabla \cdot G = 0$ (divergenceless), $\Lambda$ is a harmonic function (it satisfies Laplace’s equation):

$$\nabla \cdot G(t, x) = 0 \quad \Rightarrow \quad \nabla \cdot \nabla \Lambda(t, x) = \nabla^2 \Lambda(t, x) = 0.$$ 

For boundary conditions at infinity, this implies that $\Lambda(t, x) = \text{constant}$. For restricted regions of space (and therefore local boundary conditions), $\Lambda$ may be non-trivial.]

However, in general, a significant gauge transformation involves changing both $A$ and $\Phi$. We can generate another Lorenz-family gauge with any transformation of the following form:

**Given that**

$$\left( \nabla^2 - \frac{1}{c^2} \frac{\partial^2}{\partial t^2} \right) \Lambda(t, x) = \Box \Lambda(t, x) = 0, \quad \text{then}$$

$$A'(t, x) = A(t, x) + \nabla \Lambda(t, x) \quad \Phi'(t, x) = \Phi(t, x) - \frac{\partial \Lambda(t, x)}{\partial t}$$

[Jac 6.19 p241]

We can see that the new fields are also Lorenz gauge by simply plugging into the gauge definition above:

$$\nabla \cdot A' + \frac{1}{c^2} \frac{\partial \Phi'}{\partial t} = \nabla \cdot \left( A + \nabla \Lambda \right) + \frac{1}{c^2} \frac{\partial}{\partial t} \left( \Phi - \frac{\partial \Lambda}{\partial t} \right) = \nabla \cdot A + \nabla^2 \Lambda + \frac{1}{c^2} \frac{\partial \Phi}{\partial t} - \frac{1}{c^2} \frac{\partial^2 \Lambda}{\partial t^2}$$

$$= \left( \nabla \cdot A + \frac{1}{c^2} \frac{\partial \Phi}{\partial t} \right) + \left( \nabla^2 \Lambda - \frac{1}{c^2} \frac{\partial^2 \Lambda}{\partial t^2} \right) = 0$$

This is much easier to see relativistically, where the constraint on $\Lambda$ is:

$$\partial_\mu \partial^\mu \Lambda(t, x) = 0. \quad \text{Then}$$

$$A'^\mu = A^\mu + \partial^\mu \Lambda \quad \text{and} \quad \partial_\mu A'^\mu = \partial_\mu A^\mu + \partial_\mu \partial^\mu \Lambda = 0$$

Lorenz gauges are Lorentz covariant: the 4-divergence $\partial_\mu A^\mu = 0$. The Lorenz gauge makes the wave equations nice and symmetric, and has a simple, covariant relativistic form [Jac p241].

**Transverse, Radiation, or [gag] “Coulomb” Gauges**

Another family of gauges is radiation gauge, aka transverse gauge, or (misleadingly) Coulomb gauge, defined by
\[ \nabla \cdot \mathbf{A}(t, \mathbf{x}) = 0 \quad \text{radiation (aka transverse) gauge} \] 

which implies [Jac p241]:

\[ \Rightarrow \nabla^2 \Phi(t, \mathbf{x}) = -4\pi \rho(t, \mathbf{x}) \quad \text{(Gaussian units)} \] 

[Jac p242b] says that Coulomb gauge “indicates” that Coulomb forces propagate instantly. This is misleading, since \( \Phi \) is a mathematical field, not a physical one. However, [Jac] is quite clear that all EM fields propagate at the finite speed of light, and all behavior is consistent with Special Relativity.

We can generate another radiation-family gauge by adding a static gradient of a harmonic function to \( \mathbf{A} \):

\[ \mathbf{A}(t, \mathbf{x}) \rightarrow \mathbf{A}(t, \mathbf{x}) + \nabla \Lambda(\mathbf{x}) \quad \text{where} \quad \nabla^2 \Lambda(\mathbf{x}) = 0 \text{ preserves radiation gauge}. \]

Since \( \nabla \Lambda \) is irrotational, it preserves \( \mathbf{B} \). Since \( \partial (\nabla \Lambda) / \partial t = 0 \), it preserves \( \mathbf{E} \). [As before, for all of space, \( \Lambda \) being harmonic implies that \( \Lambda(\mathbf{x}) = \text{constant} \). For restricted regions of space, \( \Lambda \) may be non-trivial.]

The radiation gauge is often used when no significant sources are nearby, such as the radiation fields far from the source. Then \( \mathbf{E} \rightarrow 0 \) in the radiation gauge, and \( \mathbf{A} \) satisfies the homogeneous (i.e. no-source) wave-equation:

\[ \nabla^2 \mathbf{A}(t, \mathbf{x}) - \frac{1}{c^2} \frac{\partial^2 \mathbf{A}}{\partial t^2} = 0 \quad \text{[from Jac 6.30 p242]} \quad \mathbf{E}(t, \mathbf{x}) = -\frac{\partial \mathbf{A}}{\partial t}, \quad \mathbf{B}(t, \mathbf{x}) = \nabla \times \mathbf{A} \quad \text{[Jac 6.31 p242]}.
\]

With this additional condition that \( \Phi = 0 \), the \( \mathbf{E} \)-field is given by \( \mathbf{E} = -(\partial \mathbf{A}/\partial t) \). Note that:

\[ \text{Far from sources, when } \Phi = 0, \text{ the radiation gauge is also a Lorenz gauge.} \]

This follows immediately from the definitions:

\[ \nabla \cdot \mathbf{A}(t, \mathbf{x}) = 0, \quad \Phi(t, \mathbf{x}) = 0 \quad \Rightarrow \quad \nabla \cdot \mathbf{A}(t, \mathbf{x}) + \frac{1}{c^2} \frac{\partial \Phi(t, \mathbf{x})}{\partial t} = 0. \]

Furthermore, the Lorenz gauge is Lorentz covariant, but neither of the radiation gauge conditions \( \Phi = 0 \), nor \( \nabla \cdot \mathbf{A} = 0 \) is covariant (i.e., they are frame dependent).

Even a static \( \mathbf{E} \)-field can use a gauge in which \( \Phi = 0 \). This implies that \( \mathbf{A} \) contains an irrotational part which grows unboundedly (linearly) with time. (This part has no curl, and so contributes nothing to the \( \mathbf{B} \) field.) Such an unbounded \( \mathbf{A} \)-field, in this gauge, makes it hard to believe the \( \mathbf{A} \)-field is a physical field [but see the Bohm-Aharonov effect in quantum mechanics].

In general, for dynamic fields, the scalar potential \( \Phi(t, \mathbf{x}) \) is not the electric potential. In fact, in general, no time-varying “electric potential” is possible (why not?? Note that such a field is usually not conservative). The radiation gauge is no exception. \( \Phi(t, \mathbf{x}) \) is a mathematical convenience, whose relation to \( \rho(t, \mathbf{x}) \) looks the same as a static field, but whose meaning is very different. An electric potential cannot be constructed, and is certainly not simply the integral of the static Coulomb potential over the charge distribution [ref ??].

### Canonical Momentum Is Gauge Dependent

Gauge choices often help tremendously in revealing physics properties. Many calculations are greatly simplified by a good gauge choice, which makes it much easier to understand some phenomena. For example, in quantum field theory, there is one property that is very clear in one gauge, and very obscure in a 2nd gauge. But there is a 2nd property that is clear in the 2nd gauge, but obscure in the first. However,
because we know the two results are related by a gauge transformation, we know that both properties must always hold. It is important to be clear, though, which quantities are gauge invariant, and which are gauge dependent. Canonical momentum is gauge dependent.

Recall that the EM Lagrangian for a charged particle, and its canonical momentum are (gaussian units):

\[ L(q, \dot{q}) = T(q) - V(q) + \frac{e}{c} \cdot \mathbf{A}(q) \quad \text{or} \quad L = T(q_1, \ldots, q_n) - V(q_1, \ldots, q_n) + \frac{e}{c} \sum_{j=1}^{n} \dot{q}_j A_j(q_1, \ldots, q_n) \]

where \( q = (q_1, \ldots, q_n) \) are the set of generalized coordinates for the particle

\[ \mathbf{p} = \frac{\partial L}{\partial \dot{q}} = m \dot{q} + \frac{e}{c} \mathbf{A} \quad \text{or} \quad p_i = m \dot{q}_i + \frac{e}{c} A_i \]

where \( e \) is the charge of a particle.

Because of gauge freedom in choosing \( \mathbf{A}(q) \), the canonical momentum is gauge dependent, and even which components of momentum are conserved is gauge dependent. Consider a positive charge moving in a uniform \( \mathbf{B} \)-field directed out of the page. Here are 2 different \( \mathbf{A} \) fields that describe this situation:

\[ \mathbf{A}(x, y) = x \mathbf{j} \]

\[ \mathbf{A}(x, y) = -y \mathbf{i} \]

\[ \text{Figure 5.3} \quad \text{Two different gauges for } \mathbf{A} \text{ that describe the same physical situation. On the left, } p_y \text{ is conserved and } p_x \text{ is not. On the right, } p_x \text{ is conserved and } p_y \text{ is not.} \]

Recall from the Euler-Lagrange equations of motion for any Lagrangian:

\[ \frac{d}{dt} \frac{\partial L(q_i, \dot{q}_i, t)}{\partial \dot{q}_j} = \frac{\partial L}{\partial q_j} \quad \Rightarrow \quad \text{If } \frac{\partial L}{\partial \dot{q}_j} = 0, \text{ then } \frac{\partial L}{\partial q_j} = 0 = \frac{d}{dt} \frac{\partial L}{\partial \dot{q}_j} = \frac{d}{dt} p_j \]

In the situation above, the particle motion is a circle. In the left case,

\[ \mathbf{A}(x, y) = x \mathbf{j}, \quad L(x, y) = \frac{1}{2} m \dot{v}^2 + \mathbf{v} \cdot \mathbf{A} = \frac{1}{2} m \left( \dot{x}^2 + \dot{y}^2 \right) + \dot{y} x \]

\[ \frac{\partial L(x, y)}{\partial y} = 0 \quad \Rightarrow \quad p_y = \text{const} = \frac{\partial L}{\partial \dot{y}} = m \dot{y} + \frac{e}{c} A_y = m \dot{y} + \frac{e}{c} x \]

and \( p_x \) is not conserved. You can see this on the left and right edges of the circle of motion. On the left side, \( \mathbf{v} \) and \( \mathbf{A} \) point up, and add. On the right, \( \mathbf{v} \) opposes \( \mathbf{A} \), and they subtract. The sum in both cases is the same, i.e. \( p_x \) is conserved. But in the right case,

\[ \mathbf{A}(x, y) = -y \mathbf{i}, \quad L(x, y) = \frac{1}{2} m \left( \dot{x}^2 + \dot{y}^2 \right) - \dot{x} y \]

\[ \frac{\partial L}{\partial x} = 0 \quad \Rightarrow \quad p_x = \text{const} = \frac{\partial L}{\partial \dot{x}} = m \dot{x} + \frac{e}{c} A_x = m \dot{x} - \frac{e}{c} y \]

and \( p_y \) is not conserved. Here, the top and bottom edges of the circle are examples of \( \mathbf{v} \) and \( \mathbf{A} \) combining to the same \( p_x \) value at both points.
Note that any rotation of the $\mathbf{A}$ field would produce exactly the same physics, and would be just another gauge change. If we made the $\mathbf{A}$ field point at $45^\circ$, then the linear combination $p_x + p_y$ would be conserved, but neither one separately would be conserved.

**Reflection Symmetry**

When considering magnetic systems, one might encounter reflection symmetry, which is essentially a parity transformation. In such cases, we must distinguish between “polar” vectors and “axial” vectors. In short, **polar vectors** flip direction under a parity transformation, and **axial vectors** do not. So position, momentum, and E-fields, for example, are polar vectors. Angular momentum and B-fields are axial vectors (some axial vectors point along an axis of rotation, e.g. angular momentum, hence the name “axial”). This means that when you consider “reflection” symmetry (i.e., parity transformation) in E&M, you have to flip the E-field, but not the B-field. This is a little weird. You have to be careful with reflection symmetry and magnetic fields.

**Bandwidth**

[Taken from my original “Simple English Wikipedia: Bandwidth” page. As such, this section, and only this section, is public domain.] Bandwidth is a measure of how much frequency space of a spectrum is used. To clarify this, we must define some terms. Many systems work by means of vibrations, or oscillations. Vibrations are something that goes back and forth at a regular rate. Each complete cycle of “back and forth” is called, simply enough, a **cycle**. The number of cycles per second of a system is its **frequency**.

Frequency is measured in cycles per second, usually called “Hertz”, and abbreviated “Hz”.

However, most systems don’t operate at just a single frequency. They operate at many different frequencies. For example, sound is vibrations. Therefore, it has at least one frequency, and usually many different frequencies. People can hear sound frequencies as low as about 20 Hz, and as high as about 20,000 Hz. A **band** of frequencies is a continuous range of frequencies; in this example, the band of frequencies people can hear is from 20 Hz to 20,000 Hz.

Finally, bandwidth is how wide the frequency band is, that is, the highest frequency minus the lowest frequency. In the hearing example, the bandwidth of a person’s ears is about 20,000 Hz - 20 Hz = 19,980 Hz.

Bandwidth is often applied to the electromagnetic spectrum: radio waves, light waves, X-rays, and so on. Radio waves are oscillations of electric and magnetic fields. For example, the lowest United States AM radio channel covers the band of frequencies from 535,000 Hz to 545,000 Hz. It therefore has a bandwidth of 10,000 Hz (10 kHz). All United States AM radio stations have a bandwidth of 10,000 Hz. The lowest United States FM radio channel covers the band from 88,000,000 Hz (88 MHz) to 88,200,000 Hz (88.2 MHz). It therefore has a bandwidth of 200,000 Hz (200 kHz).

The term “bandwidth” has been misappropriated into the field of digital data communication. It is often incorrectly used to mean “data carrying capacity”. However, there is no such thing as “digital bandwidth”. The proper term for the data carrying capacity of a communication channel is **channel capacity**.

It is true that, in general, the channel capacity of a system increases with the bandwidth used for communication. However, many other factors come into play. As a result, in many (if not most) real systems, the channel capacity is not easily related to the channel bandwidth.

**Future Funky Electromagnetic Topics**

Let me know what topics you’d like to see.
6 Appendices

References


Griffiths

Helliwell, *Special Relativity*, ??


Jackson, 3rd ed.

Marion & Thornton, 4th ed.

Glossary

Definitions of common Electro-magnetic terms:

\(<x>\) the average (sometimes called “expectation”) value of ‘x’

**current** In a wire, current is flow rate of charge *passing a point*: \(i = \frac{dq}{dt}\) In contrast to voltage, which is *between two points*.

**fact** A small piece of information backed by solid evidence (in hard science, usually repeatable evidence). If someone disputes a fact, it is still a fact. “If a thousand people say a foolish thing, it is still a foolish thing.” See also “speculation,” and “theory.”

**linear** scales with multiplication by a constant, and summing of two parts. In other words, it *commutes with scalar multiplication, and distributes over addition*. E.g., a linear function satisfies \(f(ka + b) = kf(a) + f(b)\).

**sense** Every vector lies along an axis. The sense of a vector is simply which way the vector points along the axis. In some cases, we know the axis of a vector, but not its sense.

**speculation** A guess, possibly hinted at by evidence, but not well supported. Every scientific fact and theory started as a speculation. See also “fact,” and “theory.”

**theory** the highest level of scientific achievement: a quantitative, predictive, testable model which unifies and relates a body of facts. A theory becomes accepted science only after being supported by overwhelming evidence. A theory is not a speculation, e.g. Maxwell’s electromagnetic theory. See also “fact,” and “speculation.”

**voltage** Electrostatic potential energy difference per unit charge between two points. E.g., if \(V_{AB} = +2\) V, it means point A has 2 joules per coulomb higher potential energy than point B. If I move 3 C of charge from B to A, I must do 6 J of work on them, which is stored as 6 J of potential energy. If I let the charges run free from A to B, the E-field does 6 J of work on them, so they acquire 6 J of energy (in this, kinetic energy).

**WLOG or WOLOG** without loss of generality
Formulas

Basics

<table>
<thead>
<tr>
<th>SI</th>
<th>Gaussian</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \nabla \cdot \mathbf{D} = \rho )</td>
<td>( \nabla \cdot \mathbf{D} = 4\pi \rho )</td>
</tr>
<tr>
<td>( \mathbf{E} = \mathbf{D} - \mathbf{P} = \frac{\mathbf{D}}{\varepsilon} )</td>
<td>( \mathbf{E} = \mathbf{D} - 4\pi \mathbf{P} = \frac{\mathbf{D}}{\varepsilon} )</td>
</tr>
<tr>
<td>( \nabla \times \mathbf{E} = -\frac{\partial \mathbf{B}}{\partial t} )</td>
<td>( \nabla \times \mathbf{E} = -\frac{1}{c} \frac{\partial \mathbf{B}}{\partial t} )</td>
</tr>
<tr>
<td>( \mathbf{B} = \mathbf{H} + \mathbf{M} = \mu \mathbf{H} )</td>
<td>( \mathbf{B} = \mathbf{H} + 4\pi \mathbf{M} = \mu \mathbf{H} )</td>
</tr>
<tr>
<td>( \nabla \cdot \mathbf{B} = 0 )</td>
<td>( \nabla \cdot \mathbf{B} = 0 )</td>
</tr>
<tr>
<td>( \mathbf{F} = q \mathbf{v} \times \mathbf{B} )</td>
<td>( \mathbf{F} = q \frac{\mathbf{v} \times \mathbf{B}}{c} )</td>
</tr>
<tr>
<td>( \nabla \times \mathbf{H} = \mathbf{J} + \frac{\partial \mathbf{D}}{\partial t} )</td>
<td>( \nabla \times \mathbf{H} = \frac{4\pi}{c} \mathbf{J} + \frac{1}{c} \frac{\partial \mathbf{D}}{\partial t} )</td>
</tr>
</tbody>
</table>

\[ A(r) = \frac{\mu_0}{4\pi} \int \frac{J(r')}{|r-r'|} \, d^3 r' \quad [\text{Jac 5.32 p181}] \]

\[ A(r) = \int \frac{J(r')}{|r-r'|} \, dr' \quad [\text{LL 43.5 p111}] \]

Vector Stuff

\[ \nabla \Phi = \mathbf{e}_r \frac{\partial \Phi}{\partial r} + \mathbf{e}_\phi \frac{1}{r \sin \theta} \frac{\partial \Phi}{\partial \phi} + \mathbf{e}_\theta \frac{1}{r} \frac{\partial \Phi}{\partial \theta} \]

\[ \nabla \cdot \mathbf{A} = \frac{1}{\rho} \frac{\partial}{\partial \rho} \rho A_\rho + \frac{1}{r \sin \theta} \frac{\partial}{\partial \theta} (r A_\theta) + \frac{1}{r} \frac{\partial}{\partial \phi} A_\phi \]

\[ \nabla \times \mathbf{A} = \mathbf{e}_r \left( \frac{\partial A_\phi}{\partial \theta} - \frac{\partial A_\theta}{\partial \phi} \right) + \mathbf{e}_\phi \left( \frac{\partial A_\theta}{\partial z} - \frac{\partial A_z}{\partial \theta} \right) + \mathbf{e}_\theta \left( \frac{\partial A_z}{\partial r} - \frac{\partial A_r}{\partial z} \right) \]

\[ \nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2} \]

\[ \nabla \times (\nabla \times \mathbf{a}) = \nabla (\nabla \cdot \mathbf{a}) - \nabla^2 \mathbf{a}, \quad \mathbf{a} \times (\mathbf{b} \times \mathbf{c}) = \mathbf{b} (\mathbf{a} \cdot \mathbf{c}) - \mathbf{c} (\mathbf{a} \cdot \mathbf{b}) \]
Solutions to Laplace’s Equations

2D rectangular: \[ \Phi(x, y) = Ax + By + D \]

or \[ \Phi(x, y) = X(x)Y(y) \quad \nabla^2 \Phi = 0 \quad \Rightarrow \quad X'' + \frac{1}{X(x)} X'' + \frac{1}{Y(y)} Y'' = 0 \]

\[
\begin{cases}
X(x) = A \sin \alpha x + B \cos \alpha x, & \text{far-field } \rightarrow \text{ lowest } \alpha \\
Y(y) = C \sinh \alpha y + D \cosh (\alpha x) & \quad \text{or} \quad Y(y) = C \exp (\alpha y) + D \exp (-\alpha y)
\end{cases}
\]

or \[ \begin{cases}
X(x) = A \sin \alpha y + B \cosh (\alpha x) & \quad \text{or} \quad X(x) = A \exp (\alpha y) + B \exp (-\alpha y) \\
Y(y) = C \sin \alpha x + D \cos \alpha x, & \text{far-field } \rightarrow \text{ lowest } \alpha
\end{cases}
\]

2D polar: \[ \Phi(r, \phi) = B_0 \ln \frac{L}{A_0} + \sum_{\nu = 1}^{\infty} \left( A_{\nu} r^\nu + B_{\nu} r^{-\nu} \right) \left( C_{\nu} \sin \nu \phi + D_{\nu} \cos \nu \phi \right) \]

3D cylindrical: \[ \Phi(r, \phi, z) = R(r)Q(\phi)Z(z) \]

\[ R(r) = CJ_m (kr) + DN_m (kr) \]
\[ Q(\phi) = e^{im\phi} = A \sin m\phi + B \cos m\phi \]
\[ Z(z) = C \sinh \alpha y + D \cosh (\alpha x) \quad \text{or} \quad Z(z) = C \exp (\alpha y) + D \exp (-\alpha y) \]

\[ r = 0 \Rightarrow D = 0, \quad \Phi(r = a, \phi) = 0 \quad \Rightarrow \quad k_{mn} = \frac{x_{mn}}{a} \]

\[ f(r) = \sum_{n=1}^{\infty} A_{\nu} J_\nu \left( \frac{x_{\nu}}{a} \right) \quad A_{\nu} = \frac{2}{a^2 J_{\nu+1} (x_{\nu})} \int_0^a dr r J_\nu \left( \frac{x_{\nu}}{a} \right) \]

3D rectangular: \[ \Phi(x, y, z) = Ax + By + Cz + D \]

or \[ \Phi(x, y, z) = X(x)Y(y)Z(z) \quad \nabla^2 \Phi = 0 \quad \Rightarrow \quad \frac{1}{X(x)} X'' = \frac{1}{Y(y)} Y'' = \frac{1}{Z(z)} Z'' = 0 \]

\[ X(x) = A \sin \alpha x + B \cos \alpha x \quad \Rightarrow \quad X'' = -\alpha^2 \]

\[ Y(y) = C \sin \beta x + D \cos \beta x \quad \Rightarrow \quad Y'' = -\beta^2 \quad \alpha^2 + \beta^2 = \gamma^2, \quad \text{far-field } \rightarrow \text{ lowest } \gamma \]

\[ Z = \exp (\gamma z) \quad \text{or} \quad \sinh / \cosh (\gamma z) \quad \Rightarrow \quad Z'' = +\gamma^2 \]

\[ f(x) = \frac{A_0}{2} + A_m \cos (x) + B_m \sin (x) \quad \left\{ \begin{array}{l}
A_m \\
B_m
\end{array} \right\} = \frac{2}{a} \int_{-a/2}^{a/2} dx f(x) \left\{ \begin{array}{l}
\cos (\frac{2\pi mx}{a}) \\
\sin (\frac{2\pi mx}{a})
\end{array} \right\} \]


3D spherical: \( \Phi(r, \theta, \phi) = R(r)Q(\phi)P(\theta) \Rightarrow \)

\[
U(r) = R(r) = Ar^l + Br^{-(l+1)}, \quad Q(\phi) = e^{im\phi}
\]

\[
U_1(r) = \sqrt{\frac{2l+1}{2}} P_l(r) \quad P_0(x) = 1
\]

\[
P_l(x) = \begin{cases} 
    1 & \text{for } l = 0 \\
    \frac{1}{2} \left(3x^2 - 1\right) & \text{for } l = 1 \\
    \frac{1}{2} \left(5x^3 - 3x\right) & \text{for } l = 2 \\
    \frac{1}{8} \left(35x^4 - 30x^2 + 3\right) & \text{for } l = 3
\end{cases}
\]

Axial symmetry: \( \Phi(r, \theta) = \sum_{l=0}^{\infty} \left( A_l r^l + B_l r^{-(l+1)} \right) P_l(\cos \theta) \quad A_l = \frac{2l+1}{2} \int_0^\pi V(\theta) P_l(\cos \theta) \sin \theta \, d\theta \)

\[
Y_{lm}(\theta, \phi) = \sqrt{\frac{2l+1}{4\pi} \frac{\sin \theta}{(l+m)!}} \sum_{m=-l}^{l} \left( A_m r^l + B_m r^{-(l+1)} \right) P_l(\cos \theta) e^{im\phi} 
\]

\[
Y_{00} = \frac{1}{\sqrt{4\pi}}, \quad Y_{11} = -\frac{3}{8\pi} \sin \theta e^{i\phi}, \quad Y_{10} = \frac{3}{4\pi} \cos \theta, \quad Y_{22} = \frac{15}{4\pi} \sin^2 \theta e^{2i\phi}, \quad Y_{21} = -\frac{15}{8\pi} \sin \theta \cos \theta e^{i\phi}, \quad Y_{20} = \frac{5}{4\pi} \left(3 \cos^2 \theta - \frac{1}{2}\right)
\]

\[
g(\theta, \phi) = \sum_{l=0}^{\infty} \sum_{m=-l}^{l} A_{lm} Y_{lm}(\theta, \phi) \quad A_{lm} = \int_{\text{sphere}} d\Omega Y_{lm}^*(\theta, \phi) g(\theta, \phi)
\]

\[
\Phi(r, \theta, \phi) = \sum_{l=0}^{\infty} \sum_{m=-l}^{l} \left( A_{lm} r^l + B_{lm} r^{-(l+1)} \right) Y_{lm}(\theta, \phi) \quad \frac{1}{|x-x'|} = \sum_{l=0}^{\infty} \sum_{m=-l}^{l} r_l^l P_l(\cos \theta)
\]

Relativity and Mechanics

\[
d\tau = \frac{dx}{c}, \quad t' = \gamma \left( t - vx/c^2 \right), \quad x' = \gamma \left( x - vt \right) \quad \text{[Bra 1.55-56 p 44]}
\]

\[
\frac{1}{\gamma^2} + \beta^2 = 1 \quad u' = \frac{dx'}{d\tau} = \gamma(c, v) \quad \frac{dt}{dt'} = \frac{1 + \frac{u'v}{c^2}}{\sqrt{1 - v^2/c^2}} = \gamma \left( 1 + \frac{u'v}{c^2} \right)
\]

\[
x' = Lx = \begin{bmatrix}
    \gamma & -\gamma \beta_1 & -\gamma \beta_2 & -\gamma \beta_3 \\
    -\gamma \beta_1 & 1 + \frac{(\gamma - 1)\beta_1^2}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_1 \beta_2}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_1 \beta_3}{\beta^2} \\
    -\gamma \beta_2 & 1 + \frac{(\gamma - 1)\beta_1 \beta_2}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_2^2}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_2 \beta_3}{\beta^2} \\
    -\gamma \beta_3 & 1 + \frac{(\gamma - 1)\beta_1 \beta_3}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_2 \beta_3}{\beta^2} & 1 + \frac{(\gamma - 1)\beta_3^2}{\beta^2}
\end{bmatrix} x
\]

\[
v_{\text{total}} = \frac{v_1 + v_2}{1 + v_1 v_2 / c^2}, \quad v_y = v_y' \sqrt{1 - v^2/c^2} \quad \text{[L&L 5.1 p13]}
\]

\[
a_{\parallel} = \left( 1 - \frac{v^2}{c^2} \right)^{3/2} a' \parallel, \quad a_{\perp} = \frac{1 - v^2}{c^2} \left[ a'_{\perp} + \frac{v x (a' \times u')}{c^2} \right] \quad \text{[Jac p569]}
\]
\( E = \gamma mc^2 \quad E^2 = \left( mc^2 \right)^2 + p^2 c^2 \quad p^i = \left( E/c, p^x, p^y, p^z \right) = (E/c, p) \)

\( p = \gamma mv = Ev/c^2 \quad P = \frac{\partial L}{\partial v} = \gamma mv + \frac{q}{c} A = p + \frac{q}{c} A \)

\( \omega' = \omega (1 - \beta \cos \theta) \quad k^i = \left( \omega = ck^0, \mathbf{k} \right) \)

\( \frac{d\gamma}{dv} = \gamma^3 \frac{v}{c^2} \quad \int \frac{dv}{\gamma^3} = \int \frac{dv}{\left(1 - v^2/c^2 \right)^{3/2}} = \frac{v}{\sqrt{1 - v^2/c^2}} = \gamma v \)

\( \Omega_c = \frac{eB}{\gamma mc} = \frac{ecB}{E} \quad \frac{1}{hv'} - \frac{1}{hv} = \frac{2}{mc^2} \sin^2 \frac{\theta}{2} \)

\( L = -mc^2 \sqrt{1 - v^2/c^2} - e\phi + \frac{e}{c} \mathbf{A} \cdot \mathbf{v} \quad \left[ \text{Jac 12.12 p582} \right] \quad H = \frac{1}{2m} \left[ \left( \mathbf{P} - \frac{e}{c} \mathbf{A} \right) \right]^2 + e\phi \quad \left[ \text{LL 16.10 p49} \right] \)

\( E_\parallel' = E_\parallel \quad E_\perp' = \gamma (E_\perp + \beta \times B_\perp) \quad B_\parallel' = B_\parallel \quad B_\perp' = \gamma (B_\perp - \beta \times E_\perp) \)

Combined:

\( E' = \gamma (E + \beta \times B) - \left( \gamma - 1 \right) \beta \cdot E \quad B' = \gamma (B - \beta \times E) - \left( \gamma - 1 \right) \beta \cdot B \quad \left[ \text{Jac 11.149 p558} \right] \)

or:

\( E' = \gamma (E + \beta \times B) - \frac{\gamma^2}{\gamma + 1} \beta (\beta \cdot E) \quad B' = \gamma (B - \beta \times E) - \frac{\gamma^2}{\gamma + 1} \beta (\beta \cdot B) \quad \left[ \text{Jac 11.149 p558} \right] \)

\( v_D = c \frac{\mathbf{E} \times \mathbf{B}}{B^2} \quad v_G = \frac{\Omega_c a^2}{2B^2} (\mathbf{B} \times \mathbf{v}_\perp B) \quad v_C \approx c \frac{\gamma m}{e} \frac{v_\parallel}{R^2 B_0^2} = \frac{v_\parallel}{\Omega_c} \frac{R \times B_0}{R B_0} \quad \left[ \text{Jac 12.57-58 p590} \right] \)

\( \mu = \frac{mv_\perp}{2B} \quad \mathbf{F} = \nabla (\mu \cdot \mathbf{B}) \quad \left[ \text{Jac 5.69 p189} \right] \)

Adiabatic invariant:

\( B_a^2, \quad v_\perp^2 / B, \quad p_\perp^2 / B, \quad \gamma \mu, \quad \mu = \frac{e\Omega_c a^2}{2c} \)

NonRel:

\( \frac{d}{dt} m_\parallel = -\mu \mathbf{v} \mathbf{B} - e\mathbf{v} \nabla \phi = -\mu \frac{d\mathbf{B}}{dt} - e \frac{d\phi}{ds} = \frac{\partial}{\partial s} \left( \frac{m}{2} v_\parallel^2 \right) \quad \left[ \text{Jac 12.75 p595} \right] \)

\( \gamma \parallel \frac{\partial}{\partial s} (\gamma m_\parallel) = \frac{\partial}{\partial s} \left( \frac{m}{2} \gamma^2 v_\parallel^2 \right) \quad \gamma \approx \frac{v_\parallel}{2B_0} \frac{\partial B_\parallel}{\partial z} \quad \left[ \text{Jac 12.75 p595} \right] \)
\[ F_{ik} = \frac{\partial A_k}{\partial x^i} - \frac{\partial A_i}{\partial x^k} = \begin{bmatrix} 0 & E_x & E_y & E_z \\ -E_x & 0 & -B_z & B_y \\ -E_y & B_z & 0 & -B_x \\ -E_z & -B_y & B_x & 0 \end{bmatrix} \]

\[ F^{ik} = \begin{bmatrix} 0 & -E_x & -E_y & -E_z \\ E_x & 0 & -B_z & B_y \\ E_y & B_z & 0 & -B_x \\ E_z & -B_y & B_x & 0 \end{bmatrix} \]

\[ S = \frac{E^2 + B^2}{8\pi} \]

\[ \sigma_{\alpha \beta} = -\frac{1}{4\pi} \left( E_\alpha E_\beta + B_\alpha B_\beta - \frac{1}{2} \delta_{\alpha \beta} \left( E^2 + B^2 \right) \right) \]

\[ df = dA \cdot \hat{T} \]

\[ \mathbf{f} = \text{force} \]

\[ \mathbf{A} = \text{area} \]

\[ \mathbf{A}' = (\Phi, \mathbf{A}) \]

Faraday?

\[ \mathbf{j} = (\rho c, \mathbf{J}) \]

\[ \frac{\partial j^i}{\partial x^j} = \frac{\partial \rho}{\partial t} + \nabla \cdot \mathbf{J} = 0 \]

\[ \frac{4\pi}{c} j^i = \frac{\partial}{\partial x^i} F^{ij} \]

\[ p^i = (E/c, \mathbf{p}) \]

\[ 4\text{ force: } j^i = \frac{1}{c} F_{ik} j^k = \frac{d p^i}{d\tau} = c \frac{d}{ds} \mathbf{mu}^i = \gamma \frac{d}{dt} (\gamma mc, \gamma \mathbf{mv}) = \frac{e F^{ik} u_k}{c} \]

\[ \frac{\partial F^{ij}}{\partial x^k} + \frac{\partial F^{ki}}{\partial x^j} + \frac{\partial F^{jk}}{\partial x^i} = 0 \]

Scalars:

\[ F_{ij} F^{ij} = 2 \left( B^2 - E^2 \right) \]

\[ \epsilon^{iklm} F_{ik} F_{lm} = -4 \mathbf{E} \cdot \mathbf{B} \]

Lagrangian density:

\[ L_f = -\frac{1}{16\pi} F_{ij} F^{ij} = \frac{1}{8\pi} \left( B^2 - E^2 \right) \]

Fields:

Sheet:

\[ E_{\text{up}} = E_{\text{down}} = 2\pi \sigma \]

\[ E_{\text{conductor}} = 4\pi \sigma \]

\[ K_{\text{p}} = -c \mathbf{n} \times \mathbf{M} = c \mathbf{M} \times \mathbf{n} \]

Dipoles

\[ \mathbf{E}(\mathbf{r}) = \frac{\mathbf{P}}{r^3} \left( 2 \cos \theta \hat{\mathbf{r}} + \sin \theta \hat{\mathbf{\theta}} \right) \]

\[ \text{[Gri 3.103 p153]} \]

\[ \mathbf{M} = \frac{I}{c} \text{area} \]

\[ \Gamma = \mathbf{M} \times \mathbf{B}_e \]

\[ \mathbf{B}(\mathbf{r}) = -\nabla \frac{\mathbf{M} \cdot \mathbf{r}}{r^3} = 3 \left( \frac{M \cdot \hat{\mathbf{r}}}{r^3} \right) \hat{\mathbf{r}} - \frac{M}{cr^3} \left( 2 \cos \theta \hat{\mathbf{r}} + \sin \theta \hat{\mathbf{\theta}} \right) \]

\[ \text{[Jac 5.56 p186, Gri 3.103 p153]} \]

Energy and Work

\[ \mathbf{P} = \chi_e \mathbf{E}, \quad \mathbf{D} = \mathbf{E} + 4\pi \mathbf{P} = (1 + 4\pi \chi_e) \mathbf{E} = \varepsilon \mathbf{E} \]

\[ \mathbf{M} = \chi_m \mathbf{H}, \quad \mathbf{B} = \mathbf{H} + 4\pi \mathbf{M} = (1 + 4\pi \chi_m) \mathbf{H} = \mu \mathbf{H} \]

\[ W = \sum_{n} \phi_{n} \int_{s_n} d^2 \mathbf{s} \cdot \nabla \phi + \frac{1}{8\pi} \int d^3 r \; \rho \phi = \frac{1}{2} \sum_{n} \phi_{n} Q_{n} + \frac{1}{2} \int d^3 r \; \rho \phi = \int_{\infty} d^3 r \; \frac{E^2 + B^2}{8\pi} = \int_{\infty} d^3 r \; \frac{\nabla \phi^2 + B^2}{8\pi} \]

\[ W = \frac{1}{2} \sum_{i,j} C_{ij} \phi_i \phi_j = \frac{1}{2} C V^2 = \frac{1}{2} \frac{Q^2}{C} \quad C = \frac{2W}{V^2} = \frac{Q^2}{2W} \]

\[ w = \frac{1}{2} \mathbf{J} \cdot \mathbf{A} \quad \text{[units??]} \quad \frac{c}{\mathbf{F}} = \frac{\mathbf{J} \times \mathbf{B}}{c} \]
\[ \nabla^2 \frac{1}{|\mathbf{r}|} = \delta(0) \quad Q_{\alpha \beta} = \int_\infty d^3r' \left( 3x'_\alpha x'_\beta - \delta_{\alpha \beta} r'^2 \right) \rho(r') \quad \text{[Jac 4.9 p146, LL 41.3 p105]} \]

\[ \phi_\rho(r) = \frac{q}{\rho} + \frac{\hat{\mathbf{r}} \cdot \mathbf{d}}{\rho^2} + \frac{1}{2} \sum_{\alpha, \beta} Q_{\alpha \beta} \frac{x_\alpha x_\beta}{\rho^2} + \ldots \quad \text{[Jac 4.10 p146]} \]

\[ F = qE(0) + \nabla \left( \mathbf{P} \cdot \mathbf{E} \right)_{r=0} + \left[ \nabla \left( \frac{1}{6} \sum_{\alpha, \beta} Q_{\alpha \beta} \frac{\partial E_\alpha}{\partial x_\beta} \right) \right]_{r=0} + \ldots \quad \text{[Jac q4.5 p171]} \]

\[ \phi_\epsilon(r) = \phi_\epsilon(0) + r \cdot \nabla \phi_\epsilon(0) + \frac{1}{2} \sum_{\alpha, \beta} x_\alpha x_\beta \frac{\partial^2}{\partial x_\alpha \partial x_\beta} \phi_\epsilon(0) + \ldots \quad \text{[Jac 4.22]} \]

\[ W_\epsilon = \int_\infty d^3r \rho(r) \phi_\epsilon(r) = q \phi_\epsilon(0) + \mathbf{p} \cdot \nabla \phi_\epsilon(0) + \frac{1}{6} \sum_{\alpha, \beta} Q_{\alpha \beta} \frac{\partial^2}{\partial x_\alpha \partial x_\beta} \phi_\epsilon(0) + \ldots \]

Maxwell's Equations, Tensor

| \nabla \cdot \mathbf{D} = 4\pi \rho | \nabla \times \mathbf{E} = -\frac{1}{c} \frac{\partial \mathbf{B}}{\partial t} | \nabla \cdot \mathbf{B} = 0 | \nabla \times \mathbf{H} = \frac{4\pi}{c} \mathbf{J} + \frac{1}{c} \frac{\partial \mathbf{D}}{\partial t} |
| \iiint \mathbf{D} \cdot d\mathbf{a} = 4\pi Q | \iint \mathbf{E} \cdot d\mathbf{a} = -\frac{1}{c} \iint \frac{\partial \mathbf{B}}{\partial t} d\mathbf{a} | \iiint \mathbf{B} \cdot d\mathbf{a} = 0 | \iint \left( \frac{4\pi}{c} \mathbf{J} + \frac{1}{c} \frac{\partial \mathbf{D}}{\partial t} \right) \cdot d\mathbf{a} = \frac{4\pi}{c} \mathbf{l} |

\[ \mathbf{E} = -\nabla \phi - \frac{1}{c} \frac{\partial \mathbf{A}}{\partial t} \quad \mathbf{B} = \nabla \times \mathbf{A} = -\frac{q}{c} \frac{\mathbf{v} \times \hat{\mathbf{r}}}{r^2} = -\beta \mathbf{E}, \beta << 1 \quad d\mathbf{B} = \frac{I}{c} d\mathbf{l} \times \hat{\mathbf{r}} \quad \text{[LL 43.7]} \]

\[ \nabla^2 \mathbf{A} = -\frac{4\pi}{c} \mathbf{J} \quad \text{[LL 43.4 p110]} \quad A(r) = \frac{1}{c} \int dr' \cdot \mathbf{J}(r') \quad \text{[LL 43.5 p111]} \quad \mathbf{S} = \frac{c}{4\pi} \mathbf{E} \times \mathbf{B} \]

For \( B_0 \hat{\mathbf{z}} \):

\[ \mathbf{A} = r \frac{B_0}{2} \hat{\mathbf{\phi}} \quad \text{or} \quad \mathbf{A} = xB_0 \hat{\mathbf{x}} \quad \text{or} \quad \mathbf{A} = -yB_0 \hat{\mathbf{y}} \]

\[ \nabla \left( \nabla \times \mathbf{B} = \frac{4\pi}{c} \mathbf{J} + \frac{1}{c} \frac{\partial \mathbf{E}}{\partial t} \right) \rightarrow 0 = \frac{4\pi}{c} \left( \nabla \cdot \mathbf{J} + \frac{\partial \rho}{\partial t} \right) \quad \mathbf{F} = c \left( \frac{\mathbf{E} + \nabla \times \mathbf{B}}{c} \right) = \frac{I}{c} \mathbf{l} \times \mathbf{B} \frac{\mathbf{F}}{\mathbf{V}} = \mathbf{J} \times \mathbf{B} \]

Radiation

\[ S = |\mathbf{S}| = \frac{1}{2} \sqrt{\frac{c}{\mu} E^2} \quad \Rightarrow \quad E = \sqrt{2} \sqrt{\frac{\mu}{e}} S \quad \text{[SI, Jac 7.13 p298]} \]

\[ S = \frac{c}{4\pi} \mathbf{E} \times \mathbf{H} = \frac{c}{4\pi} E^2 \mathbf{n} \quad \Rightarrow \quad E = \sqrt{\frac{4\pi}{c} S} \quad \text{[gaussian, L&L p120]} \]
\[
\frac{d\langle P \rangle}{d\Omega} = \frac{d\sigma}{d\Omega}(l) = \frac{d\sigma}{d\Omega} c E^2
\]

\[
P = \frac{2 e^2 a^2}{3} \frac{j^2}{c^3} \quad \text{dielectric: } \frac{d\sigma}{d\Omega} = k^4 a^6 \left( \frac{e-1}{e+2} \right)^2 |\hat{E} \times \hat{r}|^2
\]

Thompson scattering:

\[
\frac{d\langle P \rangle}{d\Omega} = \frac{c}{8\pi} P^2 \sin^2 \theta \quad \frac{d\sigma}{d\Omega} = \frac{e^4}{m^2 c^4} |\hat{E} \times \hat{r}|
\]

\[
E(t) = \left[ \frac{e(1-\beta^2)(\hat{r} - \beta)}{(1-\beta \cdot \hat{r})^3} R^2 + \frac{e \hat{r} \times (\hat{r} - \beta) \times \hat{\beta}}{c (1-\beta \cdot \hat{r})^3} R \right]_{t'}
\]

\[
B(t) = [\hat{n} \times E]_{t'}
\]

\[
t = t' + \frac{r - r_0(t')}{c}
\]

\[
R(t') = r - r_0(t')
\]

Non-Rel: \[
E \approx \left[ \frac{e \hat{n} \times (\hat{n} - \beta) \times \hat{\beta}}{R} \right]_{t'} \quad \frac{dP(t)}{d\Omega} = r^2 \left( \approx R^2 \right) \hat{n} \cdot S - \frac{e^2}{4\pi c^4} |\hat{n} \times \hat{v}|^2 = \frac{e^2}{4\pi c^4} |\hat{v}|^2 \sin^2 \theta \quad \text{[Jac 14.21 p665]}
\]

\[
P(t) = \int d\Omega \frac{dP}{d\Omega} = \frac{2e^2}{3c^3} |\hat{v}|^2 \quad \text{[Jac 14.22 p665]}
\]


Multipole: \[
P(t) = \frac{2|\hat{m}|^2}{3c^3} + \frac{2|\hat{m}|^2}{3c^3} + \frac{1}{180c^3} \sum_{\alpha,\beta} [\hat{d}_{\alpha\beta}]^2 \quad \text{Relativistic: } P(t) = \frac{2e^2\gamma^4}{3c} \left[ \beta^2 - (\beta \times \hat{\beta})^2 \right]
\]

\[
\langle P \rangle = \frac{\omega^4}{3c^3} \left( |\hat{m}|^2 + |\hat{m}|^2 \right) \quad P_L(t) = \frac{2e^2\gamma^6}{3c} \beta^2 \quad P_T = \frac{2e^2}{3m^2 c^2} \gamma^4 |\hat{m}|^2 \quad \text{for fixed } \frac{dP}{dt} : P_T = \gamma^2 P_L
\]

\[
\frac{dP(t')}{d\Omega} = \frac{e^2}{4\pi c} \left[ \frac{\hat{n} \times (\hat{n} - \beta) \times \hat{\beta}}{(1-\beta \cdot \hat{n})^3} \right]_{t'} \quad \sin \theta = \hat{v} \times \hat{n}
\]

\[
\frac{dP_L(t')}{d\Omega} = \frac{e^2}{4\pi c^3} \frac{\sin^2 \theta}{(1-\beta \cos \theta)^5} \approx \frac{\theta^2}{1/2 \gamma^2 + \theta^2 / 2} \quad \text{??}
\]

\[
\frac{dP_T(t')}{d\Omega} = \frac{e^2}{2\pi c^3} \frac{|\hat{v}|^2}{(1-\beta \cos \theta)^3} \left[ 1 - \frac{\sin^2 \theta \cos^2 \phi}{\gamma^2 (1-\beta \cos \theta)^2} \right] \quad \phi = \hat{n} \text{to} \hat{\beta} \text{in} \hat{\beta} \text{??}
\]

\[\text{Reflection/Refraction}\]

\[
|E| = |B| \quad |\langle S \rangle| = \frac{c}{8\pi} \text{Re} E \times B^* = \frac{c}{8\pi} \text{Re} E \cdot E^* = \frac{c}{8\pi} |E| \quad |\langle g \rangle| = \frac{1}{8\pi c} \text{Re} E \times B^* = \frac{|E|^2}{8\pi c}
\]

\[
\hat{x} \pm i \hat{y} \rightarrow (LHCirc + helicity) / (RHCirc - helicity) \quad E_i, H_i, \text{cont}; \quad D_n, B_n \text{ cont}
\]

\[E_L : E' = \frac{2n_1 \cos \theta}{n_1 \cos \theta + (\mu_1 / \mu_2) \sqrt{n_2^2 - n_1^2 \sin^2 \theta}} \quad E'' = \frac{n_1 \cos \theta - (\mu_1 / \mu_2) \sqrt{n_2^2 - n_1^2 \sin^2 \theta}}{n_1 \cos \theta + (\mu_1 / \mu_2) \sqrt{n_2^2 - n_1^2 \sin^2 \theta}}
\]

\[E_L : E' = \frac{2n_1n_2 \cos \theta}{(\mu_1 / \mu_2)n_2 \cos \theta + n_1 \sqrt{n_2^2 - n_1^2 \sin^2 \theta}} \quad E'' = \frac{(\mu_1 / \mu_2)n_2 \cos \theta - n_1 \sqrt{n_2^2 - n_1^2 \sin^2 \theta}}{(\mu_1 / \mu_2)n_2 \cos \theta + n_1 \sqrt{n_2^2 - n_1^2 \sin^2 \theta}}
\]
\[ R = \left( \frac{S_+}{S_-} \right) = \frac{E_{+2}}{E_2} R_{\text{normal}} = \left( \frac{n_1 - n_2}{n_1 + n_2} \right)^2, \quad R_\perp = \left( \frac{\sin(\theta' - \theta)}{\sin(\theta' + \theta)} \right)^2 \]

\[ R_\parallel = \left( \frac{\tan(\theta' - \theta)}{\tan(\theta' + \theta)} \right)^2, \quad \theta + \theta' \to \frac{\pi}{2} \Rightarrow R_\parallel \to 0 \]

\[ n(\omega) = \sqrt{\varepsilon(\omega)\mu(\omega)} \quad \varepsilon_0 = \varepsilon(\omega << 1/\tau) \approx 1 + 4\pi ne^2 / m \quad \varepsilon(\omega >> \gamma) = \varepsilon_0 - \frac{4\pi nfe^2}{m\omega^2}, \]

\[ \sigma = \frac{fne^2}{m} \quad \text{decay}: \nabla^2 B = \frac{4\pi\sigma\mu}{c^2} \frac{\partial B}{\partial t} \]

\[ \delta = \sqrt{\frac{c^2}{\omega\sigma\mu}} \quad H(x) = H_0 \exp \left[ -\frac{1 - i}{\delta} x \right] \quad E(x) = \hat{z} \left( i - \frac{c}{4\pi\sigma\delta} \right) H(x) \]

\[ \frac{\text{Power}}{\text{area}} = \frac{1}{2\sigma\delta} \left( \frac{c}{4\pi} H_0 \right)^2 \quad k_{\text{eff}} = \frac{c}{4\pi} H_0 \quad \sigma_{Cu} = 10^{17} \text{ } s^{-1} ?? \]

**Waveguides**

**TM**: \( E_x(x, y) = \sin \frac{n\pi x}{a} \sin \frac{m\pi y}{b} \quad \text{TE}: \quad B_x(x, y) = \cos \frac{n\pi x}{a} \cos \frac{m\pi y}{b} \quad \gamma^2 = \left( \frac{n\pi}{a} \right)^2 + \left( \frac{m\pi}{b} \right)^2 \)

\[ \omega_{\text{min}} = \gamma_{\text{min}} c \quad \frac{\omega^2}{c^2} = k^2 + \gamma^2 \]

\[ v_{ph} = \frac{\omega}{k} = \sqrt{\frac{\gamma^2 c^2}{k^2}} \quad v_g = \frac{d\omega}{dk} = \frac{kc^2}{\sqrt{\gamma^2 c^2 + k^2 c^2}} \quad v_g v_{ph} = c^2 \]

\[ E_y = \nabla_i \phi e^{ikz-i\omega t} \quad \nabla_i^2 \phi = 0 \quad B_i = \hat{z} \times E_i \]

**Spherical Image Charges**

\[ q_{\text{out}} = -q_{\text{in}} \frac{R}{d_{\text{in}}} \quad d_{\text{out}} = \frac{R^2}{d_{\text{in}}} \quad \phi = -\lambda \ln r^2 ?? \]

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**Index**

The index is not yet developed, so go to the web page on the front cover, and text-search in this document.