

Chapter 1

Introduction

Credit

These *Mathematica* Notebooks are based on original \TeX notes by Tom R. Marsh of the Department of Physics and Astronomy at the University of Southampton.

1.1 Aims

Electromagnetism is one of the four fundamental forces. Along with gravity, it is also the one we encounter most obviously in every-day life. It is of immense practical importance and underlies optics, electricity generation, and modern communications and as well as the motors and transformers which crop up in almost every household appliance.

Electromagnetism is a *field theory*, and was the first physical theory that unified seemingly separate branches of physics, in this case optics and electricity. In field theories the physical quantities (*e.g.*, the electric and magnetic field) are defined over all space. Compare this with classical mechanics where it makes no sense to talk of the velocity of a particle defined over all space. In order to understand such continuously changing quantities we will make frequent use of vector derivatives. These are often difficult to get used to when first encountered.

The first part of the course develops these alongside *Maxwell's equations* (Chapters 2, 3), and a major aim of this course is to make you familiar with these quantities. Wave solutions of Maxwell's equations are presented as the consequences of these can be seen almost daily, even if only in a rainbow or the reflection from a puddle on a road. The aims of the course are to develop “intuition” for the behaviour of electromagnetic waves by looking at them in different situations. By the end of the course you should have become familiar with vector calculus, the physics of electric and magnetic fields, and the physics of waves, both in a general sense, and in the specific case of electromagnetic waves.

1.2 Course Structure

The combination of vector calculus and wave physics can often be difficult when first encountered. A good way to gain confidence is to apply them in problems. There will be 2 assignments followed by a tutorial where we will work through the methods of solving the problems.

1.3 Assessment

The two assignments will contribute 40% of the marks for the course and the $1\frac{1}{2}$ hour exam the remaining 60%.

1.4 Notes

The course notes are at <http://physics.uwa.edu.au/pub/Electromagnetism> as *Mathematica* Notebooks and in PDF format. The Notebooks try to be more-or-less self contained and cover everything you should know without covering too much. Please be on the look-out for errors and let me know of any that you find.

1.5 References

No one book is entirely suited to this course, and in any case books are very much a matter of personal preference. The one I like best is *Introduction to Electrodynamics* by Griffiths. *Classical Electrodynamics* by Jackson is the most famous and comprehensive text, but only recommended to the very mathematically inclined. Finally, volume 2 of *The Feynman Lectures on Physics* are worth looking at for their physical insight, particularly with regard to vector calculus. I would urge you to look at more than one treatment of any topic that you have difficulty with as each version may contain elements that help.

1. D J Griffiths, *Introduction to Electrodynamics*, 3rd edition, Prentice-Hall, 1999.
2. J D Jackson, *Classical Electrodynamics*, 2nd edition, Wiley, 1975.
3. R P Feynman, R B Leighton, and M Sands, *The Feynman Lectures on Physics, Volume 2: Electromagnetism and Matter*, Addison-Wesley, 1963-65.

1.6 Conventions

The notes are arranged in chapters each of which may cover one or more lectures. The order of the topics follows the order of the lectures. Each chapter starts with an introduction that briefly lays out what is to come. Worked examples are included, most, but not all, of which will be covered in the lectures. Some sections are marked with a \triangle : this *warning sign* indicates that you should watch out. Other sections include a \spadesuit which indicates that they contain material not covered in the lectures and not examinable. Nevertheless, they should at least be looked at in most cases. At the end of the chapters, a short section summarises the principal results and equations which you should aim to master. Appendices are used to collect together material on specific topics such as vector calculus, coordinate systems, and delta functions. The material in these appendices is examinable.

The notes follow various conventions for the symbols. Vector quantities are always in bold-face *e.g.*, \mathbf{A} . The magnitudes of vectors are scalars and are indicated by *e.g.*, A . Cross-products (wedge product) are indicated by \wedge rather than \times . Unit vectors are indicated with a hat as in $\hat{\mathbf{x}}$ for a unit vector along the x direction.

Another convention that needs to be understood is that of a *right-handed* set of axes. For many students the vector nature of electromagnetism is one of its most difficult aspects as it is often necessary to picture problems in 3-dimensions. The relative orientation of various vectors is often an issue. Starting with x and y axes at right-angles to each other, a right-handed set of axes is defined by $\hat{\mathbf{z}} = \hat{\mathbf{x}} \wedge \hat{\mathbf{y}}$. A helpful rule for cross-products is to orient your right hand so that your fingers point from the first to the second vector (*e.g.*, from $\hat{\mathbf{x}}$ to $\hat{\mathbf{y}}$ in this case). Your thumb then points in the direction of the cross-product.

Since these notes are *Mathematica* Notebooks, I use *Mathematica* conventions throughout. I find some of these conventions very useful. The exponential e (e), imaginary i (i), and differential d (d) are all displayed using "double-struck" characters (which distinguishes them from ordinary letters e , i , and d). Integrals, *e.g.*,

$$\int_0^{\frac{\pi}{2}} \cos(x) dx$$

change of variables, e.g.,

$$\frac{Dt[x]}{1+x^2} / . x \rightarrow \tan(\theta) // \text{Simplify}$$

and *total derivatives* also use d :

$$\frac{d \sin(x y)}{dx} \\ \cos(x y) \left(y + x \frac{dy}{dx} \right)$$

Partial derivatives use ∂ :

$$\frac{\partial \sin(x y)}{\partial x} \\ y \cos(x y)$$

In the figures fields and currents are indicated by crosses, \otimes , if they point down into the page and dots, \odot , if they point up out of the page.

Paul Abbott
Wednesday, July 26, 2006

Chapter 2

Gradients and Potentials

2.1 Introduction

There are many circumstances in which the rate at which a physical quantity changes with distance needs to be known. In building a road, the rate of change of height with horizontal distance — the *gradient* — is all-important. Gradients of pressure in fluids drive accelerating flows and gradients of temperature drive heat flow.

The physical quantities are usually distributed over three dimensions and so the first task in this chapter is to extend the definition of gradient from one dimension, where it is given by the derivative with respect to position, to three dimensions. We will find that the three dimensional gradient is a vector and can be calculated by application of a new operator called the *gradient* or *vector derivative* operator. We then look at how the nature of the electrostatic field allows us to define a quantity called the *potential* whose gradient is equal to the electric field.

The chapter finishes with example calculations of fields from potentials.

2.2 The Gradient

Consider a quantity such as temperature or pressure which can be assigned a value f at every point over a region. Temperature and pressure are *scalar* quantities which means that unlike vectors there is no sense in which a direction can be associated with them. Therefore f is a single number, which as it represents a physical quantity, must vary continuously. Moving from (x, y, z) to $(x + dx, y + dy, z + dz)$, the value of f changes by

$$df = \frac{\partial f}{\partial x} dx + \frac{\partial f}{\partial y} dy + \frac{\partial f}{\partial z} dz.$$

This is reminiscent of a dot product:

$$df = \left(\frac{\partial f}{\partial x} \quad \frac{\partial f}{\partial y} \quad \frac{\partial f}{\partial z} \right) \cdot \begin{pmatrix} dx \\ dy \\ dz \end{pmatrix} \equiv \left(\hat{x} \frac{\partial f}{\partial x} + \hat{y} \frac{\partial f}{\partial y} + \hat{z} \frac{\partial f}{\partial z} \right) \cdot (\hat{x} dx + \hat{y} dy + \hat{z} dz) = \nabla f \cdot d\mathbf{l}, \quad (2.1)$$

where $d\mathbf{l}$ is the *line element* vector (dx, dy, dz) , i.e.,

$$d\mathbf{l} = \begin{pmatrix} dx \\ dy \\ dz \end{pmatrix} \equiv \hat{x} dx + \hat{y} dy + \hat{z} dz,$$

and ∇ is the *vector derivative operator* (called *del* or more rarely *nabla*),

$$\nabla = \left(\frac{\partial}{\partial x} \quad \frac{\partial}{\partial y} \quad \frac{\partial}{\partial z} \right) \equiv \hat{x} \frac{\partial}{\partial x} + \hat{y} \frac{\partial}{\partial y} + \hat{z} \frac{\partial}{\partial z}$$

2.2.1 ∇f or grad f

The quantity ∇f is the *gradient* of f , and is also sometimes written as $\text{grad } f$. Since $df = \nabla f \cdot d\mathbf{l}$, for a given length of line element $d\mathbf{l}$, df is maximum when $d\mathbf{l}$ is parallel to ∇f . Thus ∇f points in the direction of maximum increase of f and its magnitude equals the rate of change of f in that direction. The gradient is the key to straightforward extension of some well-known equations that apply in one dimension. Thus the well known equation for heat conductivity:

$$Q = -\kappa \frac{dT}{dx},$$

where Q is the heat flux in $W.m^{-2}$, and κ is the conductivity, becomes

$$Q = -\kappa \nabla T \quad (2.2)$$

in 3D with the heat flux now a vector pointing in the direction of maximum decrease in temperature.

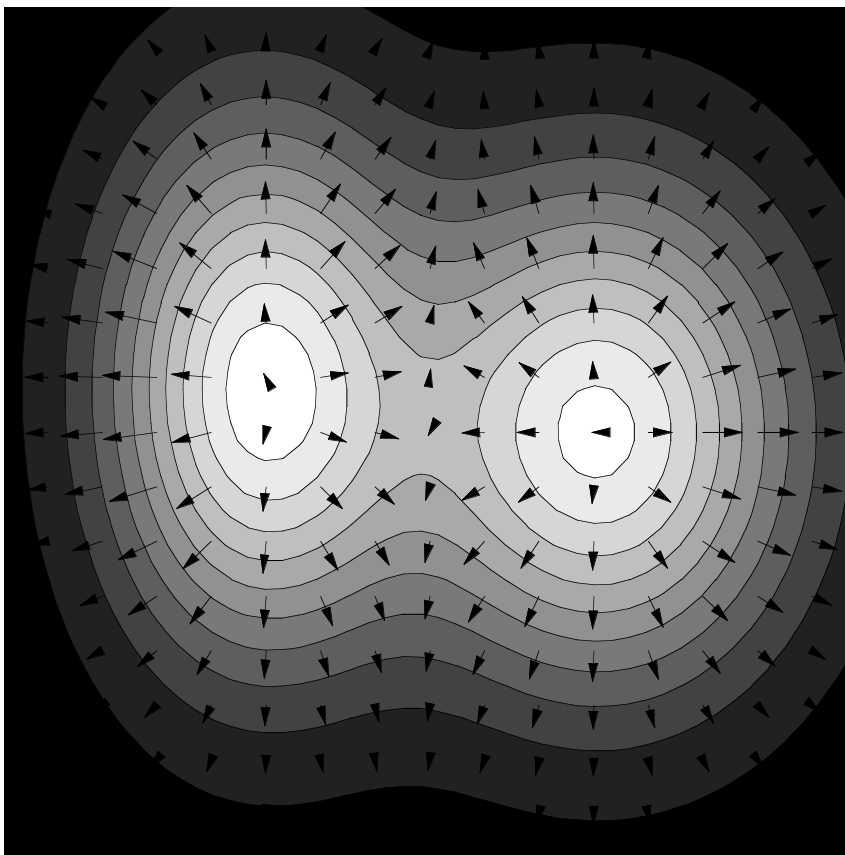


Figure 2.1 Contours of equal temperature, T , with arrows representing $-\nabla T$.

Figure 2.1 illustrates the idea of the gradient in a two-dimensional example. The contours represent lines of equal temperature (*isotherms*), in a case where there are two peaks of temperature with one higher than the other. The gradient is always *perpendicular* to the lines of equal temperature and it is large where the lines are close together.

Example 2.1 Why is the gradient always perpendicular to contour lines (or, in 3D, contour surfaces)?

If a line element $d\mathbf{l}$ lies in a line or surface along which f is constant (*i.e.*, an *isoline* or *isosurface*) then we can write $\nabla f \cdot d\mathbf{l} = 0$. Therefore $d\mathbf{l}$ must be *perpendicular* to the gradient ∇f , which is why the arrows representing the gradient in Figure 2.1 were drawn at right-angles to the contour lines.

Exercise 2.1 In Figure 2.2 what does ∇P represent?

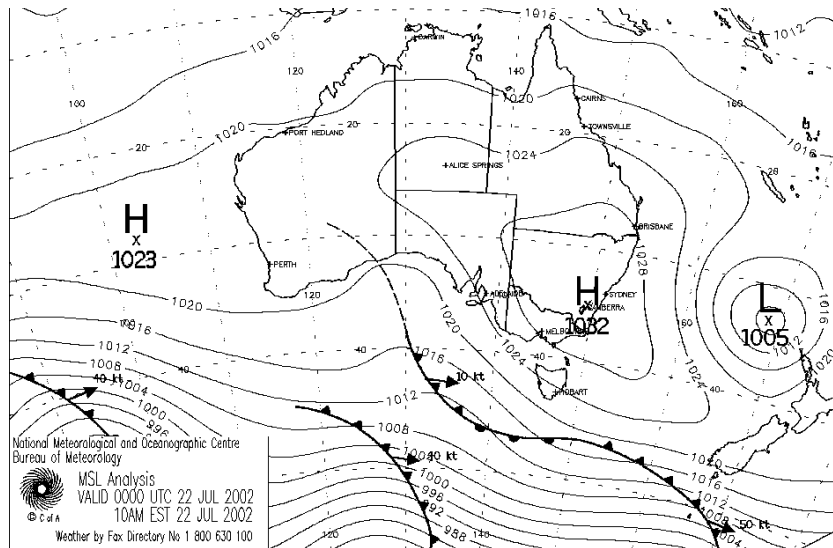


Figure 2.2 Contours of equal pressure (*isobars*), P .

2.2.2 Taylor series

You should all recall the *Taylor series* for a (differentiable) function in one variable:

$$f(x+h) = f(x) + h f'(x) + \frac{h^2}{2!} f''(x) + \frac{h^3}{3!} f^{(3)}(x) + \dots$$

In *Mathematica* you can compute Taylor series by adding an *order term* to the function:

$$f(h+x) + O(h)^5$$

$$f(x) + f'(x)h + \frac{1}{2} f''(x)h^2 + \frac{1}{6} f^{(3)}(x)h^3 + \frac{1}{24} f^{(4)}(x)h^4 + O(h^5)$$

For example, the *Maclaurin series* (*i.e.*, the Taylor series about 0) for $\tan(x)$ is

$$\tan(x) + O(x)^{10}$$

$$x + \frac{x^3}{3} + \frac{2x^5}{15} + \frac{17x^7}{315} + \frac{62x^9}{2835} + O(x^{10})$$

Since the Taylor series involves the derivatives of a function at a point, x , if $f'(x) = 0$ and $f''(x) < 0$, then

$$f(x+h) = f(x) - \alpha h^2 + \dots,$$

(where $\alpha > 0$) and x is a *local maximum* of f because, in the neighbourhood of x (*i.e.*, around $h = 0$), f decreases as we move away from x . But of course, you already knew this from high-school calculus. However, in more than one variable, the situation is more complicated.

One very interesting (formal) way of writing the Taylor series is

$$f(x+h) = e^{h \frac{\partial}{\partial x}} f(x)$$

where this formal notation is interpreted as

$$f(x+h) = e^{h \frac{\partial}{\partial x}} f(x) = \left(1 + h \frac{\partial}{\partial x} + \frac{1}{2!} \left(h \frac{\partial}{\partial x} \right)^2 + \frac{1}{3!} \left(h \frac{\partial}{\partial x} \right)^3 + \dots \right) f(x)$$

$$= f(x) + h f'(x) + \frac{h^2}{2!} f''(x) + \frac{h^3}{3!} f^{(3)}(x) + \dots$$

This idea turns out to be useful in group theory. The action of the operator $e^{h \frac{\partial}{\partial x}}$ on $f(x)$ has the effect of *translating* the function to $f(x+h)$.

Another advantage of this notation is that it is straightforward to extend it to any number of variables by replacing $h \frac{\partial}{\partial x}$ with $\mathbf{h} \cdot \nabla$:

$$\begin{aligned} f(\mathbf{x} + \mathbf{h}) &= e^{\mathbf{h} \cdot \nabla} f(\mathbf{x}) = \left(1 + \mathbf{h} \cdot \nabla + \frac{1}{2!} (\mathbf{h} \cdot \nabla)^2 + \frac{1}{3!} (\mathbf{h} \cdot \nabla)^3 + \dots \right) f(\mathbf{x}) \\ &= f(\mathbf{x}) + (\mathbf{h} \cdot \nabla) f(\mathbf{x}) + \frac{1}{2!} (\mathbf{h} \cdot \nabla)^2 f(\mathbf{x}) + \frac{1}{3!} (\mathbf{h} \cdot \nabla)^3 f(\mathbf{x}) + \dots \end{aligned}$$

Some care needs to be taken when interpreting this expression: For two variables, the second term is

$$\begin{aligned} (\mathbf{h} \cdot \nabla) f(\mathbf{x}) &= (h, k) \cdot (\partial_x, \partial_y) f(x, y) = \\ &= h f^{(1,0)}(x, y) + k f^{(0,1)}(x, y), \end{aligned}$$

and the third term is

$$\begin{aligned} \frac{1}{2!} (\mathbf{h} \cdot \nabla)^2 f(\mathbf{x}) &= \frac{1}{2} (h, k) \cdot (\partial_x, \partial_y) ((h, k) \cdot (\partial_x, \partial_y) f(x, y)) = \\ &= \frac{1}{2} f^{(2,0)}(x, y) h^2 + k f^{(1,1)}(x, y) h + \frac{1}{2} k^2 f^{(0,2)}(x, y). \end{aligned}$$

Omitting the factor of $1/2$, this can be written in matrix notation as

$$(\mathbf{h} \cdot \nabla)^2 f(\mathbf{x}) = \mathbf{h}^T H \mathbf{h} = (h \ k) \cdot \begin{pmatrix} \frac{\partial^2 f(x,y)}{\partial x \partial x} & \frac{\partial^2 f(x,y)}{\partial x \partial y} \\ \frac{\partial^2 f(x,y)}{\partial y \partial x} & \frac{\partial^2 f(x,y)}{\partial y \partial y} \end{pmatrix} \cdot \begin{pmatrix} h \\ k \end{pmatrix},$$

where (apart from a sign), H is the *Hessian* matrix.

We require an important result from linear algebra: A *symmetric* $n \times n$ matrix M is *positive definite* $\Leftrightarrow \mathbf{x}^T M \mathbf{x} > 0$ for all $\mathbf{x} \neq \mathbf{0}$ in $\mathbb{R}^n \Leftrightarrow$ all the n eigenvalues, λ_i , of M are such that each $\lambda_i > 0$. Similarly, a *negative definite* matrix has each $\lambda_i < 0$. Since the Taylor series

$$f(x+h, y+k) = f(x, y) + (h \ k) \cdot \begin{pmatrix} \frac{\partial f(x,y)}{\partial x} \\ \frac{\partial f(x,y)}{\partial y} \end{pmatrix} + \frac{1}{2} (h \ k) \cdot \begin{pmatrix} \frac{\partial^2 f(x,y)}{\partial x \partial x} & \frac{\partial^2 f(x,y)}{\partial x \partial y} \\ \frac{\partial^2 f(x,y)}{\partial y \partial x} & \frac{\partial^2 f(x,y)}{\partial y \partial y} \end{pmatrix} \cdot \begin{pmatrix} h \\ k \end{pmatrix} + \dots$$

involves the partial derivatives of f at the point, $\mathbf{x} = (x, y)$, if $\nabla f(\mathbf{x}) = \mathbf{0}$, i.e., $f^{(1,0)}(x, y) = 0 = f^{(0,1)}(x, y)$, and H is *negative definite* then (x, y) is a *local maximum* of f .

The relationship between the sign of the eigenvalues and the sign of $\mathbf{x}^T M \mathbf{x}$ results directly from the definition of the eigenvalues, λ_i and corresponding (orthonormal) eigenvectors, \mathbf{u}_i , of a *symmetric* matrix:

$$M \mathbf{u}_i = \lambda_i \mathbf{u}_i \Rightarrow \mathbf{u}_j^T M \mathbf{u}_i = \lambda_i \mathbf{u}_j^T \mathbf{u}_i = \lambda_i \delta_{i,j}.$$

If the eigenvectors *span* \mathbb{R}^n , we can express *any* vector in \mathbb{R}^n as $\mathbf{x} = \alpha_1 \mathbf{u}_1 + \dots + \alpha_n \mathbf{u}_n$ where each $\alpha_i \in \mathbb{R}$. Then

$$\mathbf{x}^T M \mathbf{x} = \left(\sum_{j=1}^n \alpha_j \mathbf{u}_j^T \right) M \left(\sum_{i=1}^n \alpha_i \mathbf{u}_i \right) = \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j \mathbf{u}_j^T M \mathbf{u}_i = \sum_{i=1}^n \sum_{j=1}^n \alpha_i \alpha_j \lambda_i \delta_{i,j} = \sum_{i=1}^n \lambda_i \alpha_i^2.$$

Since $\alpha_i^2 > 0$ and the α_i are arbitrary, $\mathbf{x}^T M \mathbf{x} > 0 \Leftrightarrow \lambda_i > 0, \forall_{i=1,2,\dots,n}$.

The eigenvectors *diagonalize* the symmetric matrix, which can be written in the form

$$M = P^T . D . P,$$

where P is the matrix of eigenvectors, $P = (\mathbf{u}_1 | \mathbf{u}_2 | \dots | \mathbf{u}_n)$, and D is the diagonal matrix with the eigenvalues $\lambda_1, \lambda_2, \dots, \lambda_n$ along the diagonal. Then

$$\mathbf{x}^T M \mathbf{x} = \mathbf{x}^T P^T D P \mathbf{x} = (P \mathbf{x})^T D (P \mathbf{x}) \Rightarrow \mathbf{y}^T D \mathbf{y} = (\lambda_1 y_1^2 + \lambda_2 y_2^2 + \dots + \lambda_n y_n^2),$$

where $\mathbf{y} = P \mathbf{x}$. Clearly, $\lambda_1 y_1^2 + \lambda_2 y_2^2 + \dots + \lambda_n y_n^2 > 0$ for *arbitrary* (real) y_i only if all $\lambda_i > 0$.

Example 2.2 Consider the symmetric 2×2 matrix $M = \begin{pmatrix} 5 & -2 \\ -2 & 8 \end{pmatrix}$. Is M positive definite?

If we compute $\mathbf{x}^T M \mathbf{x}$

$$M = \begin{pmatrix} 5 & -2 \\ -2 & 8 \end{pmatrix};$$

`{x, y}.M.{x, y} // Factor`

$$5x^2 - 4xy + 8y^2$$

it is not immediately obvious that this expression is positive for arbitrary x and y . However, if we write the result in the form

$$\% == \frac{9}{5} (x - 2y)^2 + \frac{4}{5} (2x + y)^2 // \text{Simplify}$$

True

it is now obvious, since $(x - 2y)^2$ and $(2x + y)^2$ are both positive for all $\mathbf{x} = (x, y) \neq \mathbf{0}$ in \mathbb{R}^2 . Alternatively, we see that both eigenvalues are positive:

`$\Lambda = \text{Eigenvalues}[M]$`

{9, 4}

Hence M is positive definite. Alternatively, with

`$D = \text{DiagonalMatrix}[\Lambda]$`

$$\begin{pmatrix} 9 & 0 \\ 0 & 4 \end{pmatrix}$$

then clearly $\mathbf{y}^T D \mathbf{y} = \lambda_1 u^2 + \lambda_2 v^2 > 0$ for all $\mathbf{y} = (u, v) \neq \mathbf{0}$ in \mathbb{R}^2 .

`{u, v}.D.{u, v}`

$$9u^2 + 4v^2$$

The *orthogonal* eigenvectors are

`$\text{Eigenvectors}[M]^T$`

$$\begin{pmatrix} -1 & 2 \\ 2 & 1 \end{pmatrix}$$

We need to make these *orthonormal*:

$$P = \frac{\%}{\text{Norm} /@ \%} \\ \begin{pmatrix} -\frac{1}{\sqrt{5}} & \frac{2}{\sqrt{5}} \\ \frac{2}{\sqrt{5}} & \frac{1}{\sqrt{5}} \end{pmatrix}$$

$$P^T \cdot P$$

$$\begin{pmatrix} 1 & 0 \\ 0 & 1 \end{pmatrix}$$

P diagonalizes the matrix M .

$$P^T \cdot M \cdot P \text{ // Simplify}$$

$$\begin{pmatrix} 9 & 0 \\ 0 & 4 \end{pmatrix}$$

and we confirm that $P^T D P = M$:

$$P^T \cdot \mathcal{D} \cdot P == M$$

True

Computing $(P \mathbf{x})^T D (P \mathbf{x})$ we also obtain a result that is positive for all $\mathbf{x} = (x, y) \neq \mathbf{0}$ in \mathbb{R}^2 .

$$(P \cdot \{x, y\}) \cdot \mathcal{D} \cdot (P \cdot \{x, y\})$$

$$4 \left(\frac{2x}{\sqrt{5}} + \frac{y}{\sqrt{5}} \right)^2 + 9 \left(\frac{2y}{\sqrt{5}} - \frac{x}{\sqrt{5}} \right)^2$$

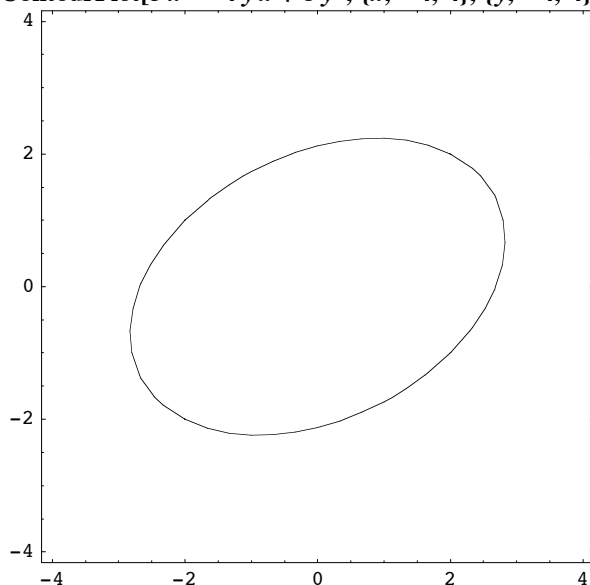
$$\text{Simplify /@ \%}$$

$$\frac{9}{5} (x - 2y)^2 + \frac{4}{5} (2x + y)^2$$

Example 2.3 Describe the conic $5x^2 - 4xy + 8y^2 = 36$.

Visualizing the conic shows that it is an ellipse:

`ContourPlot[5 x2 - 4 y x + 8 y2, {x, -4, 4}, {y, -4, 4}, Contours -> {36}, ContourShading -> False];`



Write the equation in matrix form:

$$M = \begin{pmatrix} 5 & -2 \\ -2 & 8 \end{pmatrix};$$

$$\{x, y\} \cdot M \cdot \{x, y\} == 36 \text{ // ExpandAll}$$

$$5x^2 - 4yx + 8y^2 = 36$$

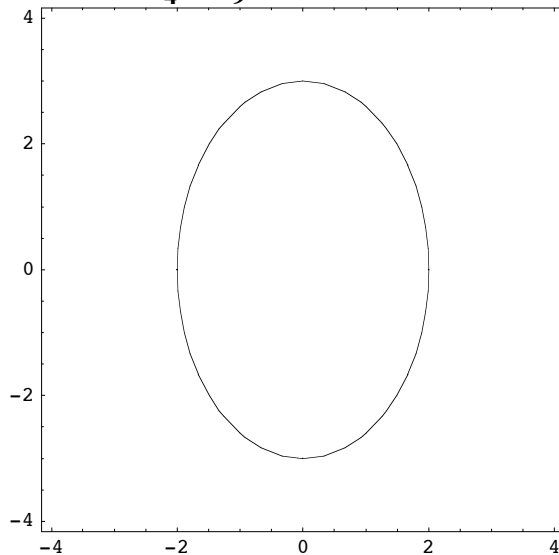
Diagonalizing the matrix, $\mathbf{x}^T M \mathbf{x} = \mathbf{x}^T P^T D P \mathbf{x} = (P \mathbf{x})^T D (P \mathbf{x}) = \mathbf{y}^T D \mathbf{y}$, where $\mathbf{y} = (u, v) = P \mathbf{x}$. Hence the equation becomes

$$\{u, v\} \cdot \mathcal{D}\{u, v\} == 36 // \text{Simplify}$$

$$9u^2 + 4v^2 == 36$$

that is, $\frac{u^2}{4} + \frac{v^2}{9} = 1$, which is the equation of an *ellipse*.

`ContourPlot` $\left[\frac{u^2}{4} + \frac{v^2}{9}, \{u, -4, 4\}, \{v, -4, 4\}, \text{Contours} \rightarrow \{1\}, \text{ContourShading} \rightarrow \text{False}\right];$



The effect of the orthogonal matrix P on x is to *rotate* the axes:

$$P \cdot \{x, y\}$$

$$\left\{ \frac{2y}{\sqrt{5}} - \frac{x}{\sqrt{5}}, \frac{2x}{\sqrt{5}} + \frac{y}{\sqrt{5}} \right\}$$

2.2.3 Conservative fields and potentials

Eq. 2.1, $df = \nabla f \cdot d\mathbf{l}$, can be used to calculate the finite change in f when moving from two points, A to B (Fig. 2.3):

$$f(B) - f(A) = \int_A^B df = \int_A^B \nabla f \cdot d\mathbf{l}.$$

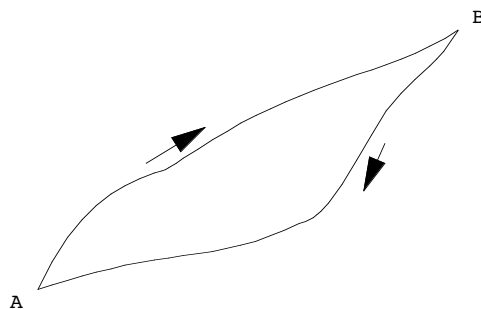


Figure 2.3 Path of integration from A to B and back again.

If we then move back from B to A over a different path, the total change in f will be zero and thus

$$\oint \nabla f \cdot d\mathbf{l} = 0,$$

where the symbol \oint indicates an integral over a *closed loop*. The reverse of this process can be shown to be true. That is, if integrals over closed loops in a vector field \mathbf{A} are always zero *i.e.*,

$$\oint \mathbf{A} \cdot d\mathbf{l} = 0$$

for *any* loop, then \mathbf{A} can be derived from a *scalar field*, called ψ say, by taking its gradient

$$\mathbf{A} = \nabla\psi.$$

This is an important theorem since it is generally much easier to work with scalars than vectors.

Since the force on a charge q in an electric field \mathbf{E} is $q\mathbf{E}$, and so a force $-q\mathbf{E}$ needs to be applied to hold the charge still, the integral

$$-\oint q\mathbf{E} \cdot d\mathbf{l}$$

represents the work needed to move the charge around a loop. In electrostatics this must be zero or else we could obtain energy indefinitely by allowing the charge to move around the loop in the direction that makes the work needed negative. In electrostatics $\oint \mathbf{E} \cdot d\mathbf{l} = 0$ for all loops and therefore from above, we must be able to derive \mathbf{E} from a scalar, *i.e.*, $\mathbf{E} = -\nabla\phi$. In fact by convention we write

$$\mathbf{E} = -\nabla\phi,$$

where ϕ is called the *electric potential*. The minus sign means that the potential increases as one nears positive charges and makes ϕ the work done in bringing a unit charge from infinity to a given point. The reasoning above breaks down in time varying cases when it is possible for $\oint \mathbf{E} \cdot d\mathbf{l} \neq 0$ (*e.g.*, think of the coils of a transformer). Thus the above equation applies in electrostatics only.

We have used the conservation of energy to argue that $\oint \mathbf{E} \cdot d\mathbf{l} = 0$ and any vector field that satisfies this condition is known as a *conservative* field. Not all fields satisfy this condition. For example any field that can be drawn in closed loops cannot have a zero-line integral around these loops. The magnetic field around a wire is one example, and in general it is not possible to derive magnetic fields from scalar potentials.

2.2.4 Calculating fields from potentials

The calculation of fields from potentials is best illustrated with some examples. We start with a simple one.

Example 2.4 What is the electric field equivalent to the potential $\phi = -kx$?

We apply the vector derivative operator through $\mathbf{E} = -\nabla\phi$ or in component form

$$\begin{pmatrix} E_x \\ E_y \\ E_z \end{pmatrix} = - \begin{pmatrix} \frac{\partial}{\partial x} \\ \frac{\partial}{\partial y} \\ \frac{\partial}{\partial z} \end{pmatrix} (-kx) = \begin{pmatrix} k \\ 0 \\ 0 \end{pmatrix}.$$

Thus $\phi = -kx$ is the potential of a uniform field pointing in the x direction (*i.e.*, \hat{i}) with magnitude k .

This example can be generalised:

Example 2.5 What is the electric field equivalent to the potential $\phi = -\mathbf{A} \cdot \mathbf{r}$ where \mathbf{A} is a constant vector and \mathbf{r} is the position vector?

The dot product can be expanded out so

$$\phi = -A_1 x - A_2 y - A_3 z.$$

We then have

$$\mathbf{E} = -\nabla\phi = \begin{pmatrix} A_1 \\ A_2 \\ A_3 \end{pmatrix} = \mathbf{A}.$$

Therefore a uniform field, \mathbf{E} , has a potential of the form $\phi = -\mathbf{E} \cdot \mathbf{r}$.

Now for a trickier case:

Example 2.6 What is the electric field equivalent to the potential $\phi = 1/r$ where r is the distance from a point?

This can be answered in two ways:

(1) The direct approach is to apply the vector derivative operator to $1/r$ remembering that $r^2 = x^2 + y^2 + z^2$. Thus since

$$\frac{\partial}{\partial x} \left(\frac{1}{r} \right) = -\frac{1}{r^2} \frac{\partial r}{\partial x} = -\frac{1}{r^2} \frac{\partial(x^2 + y^2 + z^2)^{1/2}}{\partial x},$$

and

$$\frac{\partial r}{\partial x} = \frac{\partial(x^2 + y^2 + z^2)^{1/2}}{\partial x} = \frac{1}{2} (x^2 + y^2 + z^2)^{-1/2} 2x = \frac{x}{r},$$

we obtain

$$\frac{\partial}{\partial x} \left(\frac{1}{r} \right) = -\frac{x}{r^3},$$

and so, with similar expressions for the other components, we obtain

$$\mathbf{E} = -\nabla\phi = \frac{\mathbf{r}}{r^3} \equiv \frac{\hat{\mathbf{r}}}{r^2} \equiv \frac{\mathbf{e}_r}{r^2}, \quad (2.3)$$

where $\hat{\mathbf{r}}$ or \mathbf{e}_r are unit vectors pointing in the radial direction. Therefore, as expected, a $1/r$ potential gives a $1/r^2$ electric field.

(2) A more intuitive approach can be taken based upon Eq. 2.1, $df = \nabla f \cdot d\mathbf{l}$. If $d\mathbf{l}$ is parallel to ∇f , *i.e.*, we step along the direction of the gradient, then this becomes $df = |\nabla f| dl$, or

$$|\nabla f| = \frac{df}{dl}$$

$$\phi = 1/r \Rightarrow -\frac{d\phi}{dr} = \frac{1}{r^2} \Rightarrow \mathbf{E} = \frac{\hat{\mathbf{r}}}{r^2}$$

along a path parallel to the gradient. For the electric field we can similarly write $E = -d\phi/dl$ for a path parallel to the field. For $\phi = 1/r$ the field must point in the radial direction by symmetry so we take the derivative moving out in radius, *i.e.*, $E = -d\phi/dr$ for any spherically symmetric potential ϕ . This trivially gives the result (Eq. 2.3) obtained more painfully above, and can be applied to any potential that varies with r only.

Computation of fields using Mathematica

Define the distance between two points as follows:

$$\text{Unprotect}[\text{Norm}]; \|z_ \| := \sqrt{z.z}; \text{Protect}[\text{Norm}];$$

The potential of a (point) charge q positioned at \mathbf{r}_0 , measured at \mathbf{r} is

$$\phi(\mathbf{r}_-, \mathbf{r}_0, q) := \frac{q}{\|\mathbf{r} - \mathbf{r}_0\|}$$

For example, the potential due to a unit charge at the origin

$$O = \{0, 0, 0\};$$

measured at the point

$$P = \{x, y, z\};$$

is

$$\frac{\phi(P, O)}{\sqrt{x^2 + y^2 + z^2}}$$

In *Mathematica* after defining ∇ :

$$\nabla f := \left\{ \frac{\partial f}{\partial x}, \frac{\partial f}{\partial y}, \frac{\partial f}{\partial z} \right\}$$

We can compute $-\nabla\phi$ in Cartesian coordinates directly:

$$\mathcal{E} = \text{Simplify}\left[-\nabla \frac{1}{\sqrt{x^2 + y^2 + z^2}} /. x^2 + y^2 + z^2 \rightarrow r^2, r > 0\right]$$

$$\left\{ \frac{x}{r^3}, \frac{y}{r^3}, \frac{z}{r^3} \right\}$$

Introducing the unit vector, \mathbf{e}_r , in the radial direction

$$\mathbf{e}_r = \frac{\mathbf{P}}{r}$$

$$\left\{ \frac{x}{r}, \frac{y}{r}, \frac{z}{r} \right\}$$

the electric field \mathcal{E} can be written as

$$\mathcal{E} = \frac{\mathbf{P}}{r^3} = \frac{\mathbf{e}_r}{r^2}$$

True

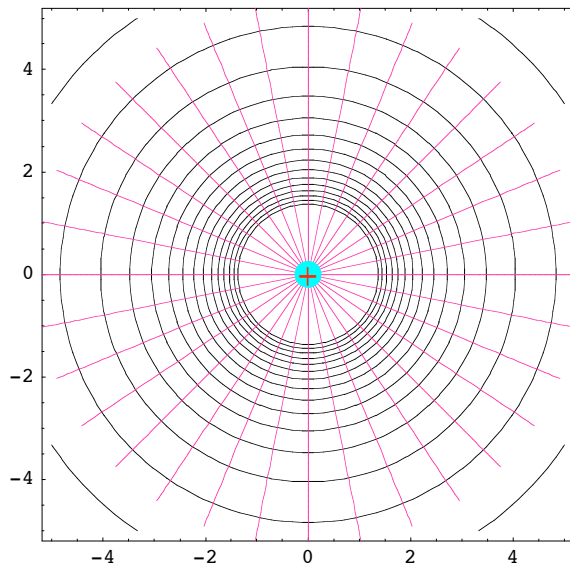
Alternatively, the electric field for any spherically symmetric potential ϕ can be computed using $E = -d\phi/dr$:

$$\mathcal{E} = -\frac{d}{dr} \frac{1}{r} \mathbf{e}_r$$

True

2.2.5 Point Charge — Monopole

The electric field and potential of a point charge (*monopole*) can be visualized as follows. Computing the *field lines* (by numerically solving a set of differential equations), we show the charge, equipotential lines (black), and field lines (purple) together:



△ Note that the field lines and equipotentials are *orthogonal* (i.e., they intersect at right-angles).

Restricting attention to the x - z plane (i.e., $y = 0$),

$$\mathbf{P} = \{x, 0, z\};$$

the potential of a point (unit) charge is

$$\phi_1(\mathbf{x}_-, \mathbf{z}_-) = \phi(\mathbf{P}, \mathbf{O}) = \frac{1}{\sqrt{x^2 + z^2}}$$

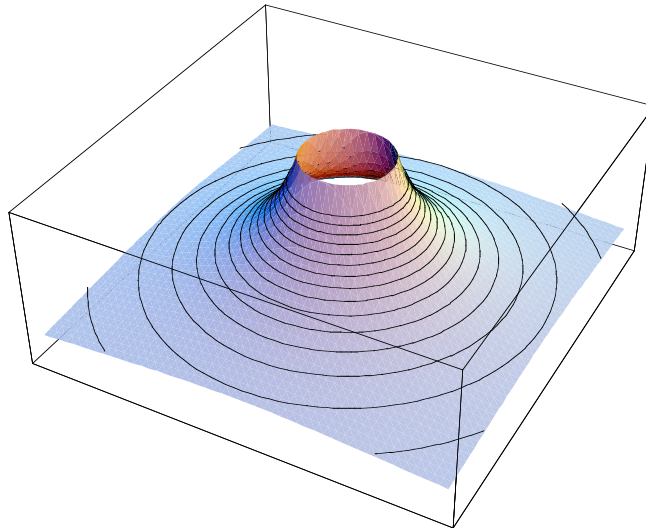
and the (Cartesian) components of the electric field are

$$\{\mathcal{E}_x, \mathcal{E}_y, \mathcal{E}_z\} = -\nabla \phi_1(x, z) = \left\{ \frac{x}{(x^2 + z^2)^{3/2}}, 0, \frac{z}{(x^2 + z^2)^{3/2}} \right\}$$

△ Note that restricting attention to the x - z plane simplifies the computations slightly and is convenient when plotting graphs of the potential and field. However, you should remember that, in general, the potential and field are functions of all 3 (Cartesian) coordinates.

△ There is an important subtlety here: from first year you should already be aware that the *density of lines* in plots of the electric field are proportional to the strength of the field. However, the density needs to be computed in 3 dimensions (i.e., lines per unit volume) rather than in 2 dimensions (i.e., lines per unit area). If you do this for a point charge you will find that the density of lines does indeed go like the inverse square of the distance from the charge, i.e., $1/r^2$.

A powerful alternative visualization is a surface plot with the equipotential lines superimposed onto the surface:



Imagine placing a ball bearing on this surface under the influence of gravity acting in the vertical direction. Qualitatively, the magnitude and direction of the force on the ball-bearing is obvious. By analogy, one can immediately obtain the forces acting on and the resulting motion of a positive test charge in such a potential.

2.2.6 Pure Dipole

The second method used in example 2.4 can be applied more generally to cases lacking symmetry and is usually the way to proceed unless the potential is given in terms of x , y , and z (in which case direct application of $\partial/\partial x$ etc, is easiest). Consider a (pure) dipole potential of the form

$$\phi = \frac{\mathbf{p} \cdot \hat{\mathbf{r}}}{4\pi\epsilon_0 r^2} = \frac{\mathbf{p} \cdot \mathbf{r}}{4\pi\epsilon_0 r^3} = \frac{p_1 x + p_2 y + p_3 z}{4\pi\epsilon_0 (x^2 + y^2 + z^2)^{3/2}} = \frac{p \cos(\theta)}{4\pi\epsilon_0 r^2} \quad (2.4)$$

where \mathbf{p} is a constant vector and θ is the angle between \mathbf{p} and the radial direction, $\hat{\mathbf{r}}$. What is the electric field of such a potential? Since the potential is expressed in spherical coordinates r and θ , it is easiest to work out the field in the radial (increasing r) and tangential (increasing θ) directions.

We start again from Eq. 2.1, $df = \nabla f \cdot d\mathbf{l}$, or its equivalent — here $d\phi = -\mathbf{E} \cdot d\mathbf{l}$. If $d\mathbf{l}$ is parallel to the radial direction, only the radial component of \mathbf{E} , E_r , contributes to the dot product and thus

$$E_r = - \left. \frac{d\phi}{dl} \right|_{\theta \text{ const}} = - \frac{dr}{dl} \frac{\partial \phi}{\partial r} = - \frac{\partial \phi}{\partial r},$$

with the *partial derivative* showing that only r changes. Similarly if we move tangentially, only the tangential component E_θ contributes to the dot product and

$$E_\theta = - \left. \frac{d\phi}{dl} \right|_{r \text{ const}} = - \frac{d\theta}{dl} \frac{\partial \phi}{\partial \theta} = - \frac{1}{r} \frac{\partial \phi}{\partial \theta}.$$

△ Note here that, for r constant, if we move from (r, θ) to $(r, \theta + d\theta)$ we have moved by $dl = r d\theta$ so $d\theta/dl = 1/r$ which is why a $1/r$ term appears (as it must to give the correct dimensions). For θ constant, if we move from (r, θ) to $(r + dr, \theta)$ we have moved by $dl = dr$ so $dr/dl = 1$.

Applying $E_r = -\frac{\partial \phi}{\partial r}$ and $E_\theta = -\frac{1}{r} \frac{\partial \phi}{\partial \theta}$ to the potential of Eq. 2.4 we find

$$E_r = \frac{2 p \cos(\theta)}{4\pi\epsilon_0 r^3},$$

$$E_\theta = \frac{p \sin(\theta)}{4\pi\epsilon_0 r^3}.$$

The total field is

$$\mathbf{E} = E_r \mathbf{e}_r + E_\theta \mathbf{e}_\theta,$$

and is illustrated in Figure 2.4 below.

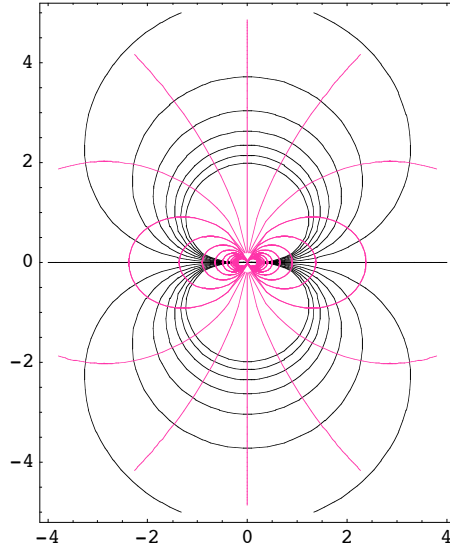


Figure 2.4 The field pattern (purple) and equipotential lines (black) for a potential of the form $p \cos(\theta)/r^2$.

Total derivative versus partial derivative

Recall the definition of the *total derivative*. For f a function of r and θ we find that

$$\frac{df(r, \theta)}{dl} = \frac{dr}{dl} f^{(1,0)}(r, \theta)$$

where $f^{(0,1)}(r, \theta)$ denotes the *partial derivative* of f with respect to its second argument, *i.e.*, θ :

$$\frac{\partial f(r, \theta)}{\partial \theta} = f^{(0,1)}(r, \theta)$$

and similarly for $f^{(1,0)}(r, \theta)$. Computing the total derivative for constant θ we obtain:

$$\text{SetAttributes}[\theta, \text{Constant}]; \frac{df(r, \theta)}{dl} = \frac{dr}{dl} f^{(1,0)}(r, \theta)$$

and similarly for constant r .

2.2.7 Dipole

The total potential, $\phi_2(x, z)$, of a pair of equal and opposite charges, $+1$ positioned at $\{0, 0, 1\}$ and -1 positioned at $\{0, 0, -1\}$, is

$$\phi_2(\underline{x}_-, z_-) = \frac{\phi(P, \{0, 0, -1\}, -1)}{1} + \frac{\phi(P, \{0, 0, 1\}, 1)}{1}$$

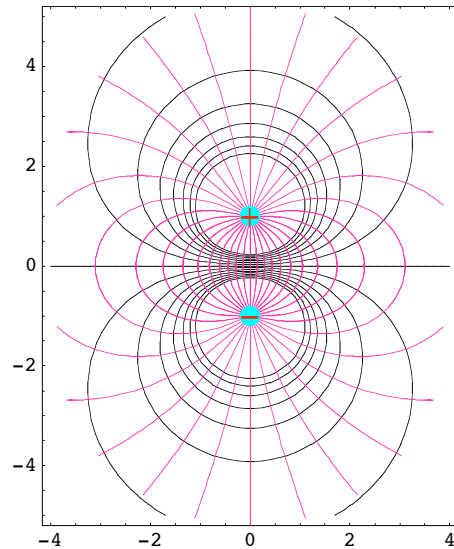
$$\frac{1}{\sqrt{x^2 + (z-1)^2}} - \frac{1}{\sqrt{x^2 + (z+1)^2}}$$

with corresponding electric field

$$\{\mathcal{E}_x, \mathcal{E}_y, \mathcal{E}_z\} = -\nabla\phi_2(x, z)$$

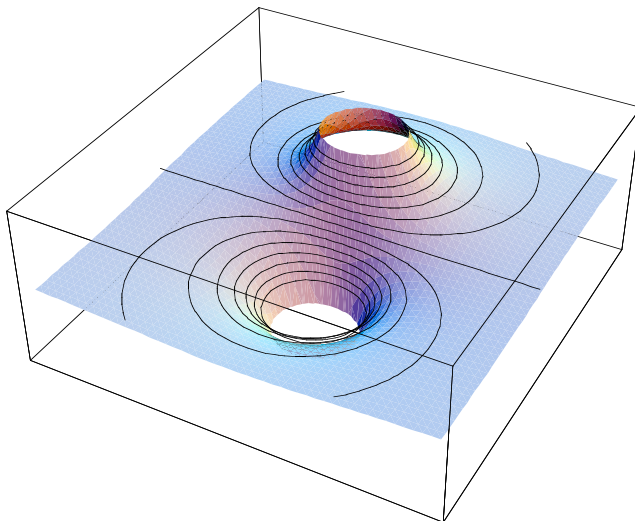
$$\left\{ \frac{x}{(x^2 + (z-1)^2)^{3/2}} - \frac{x}{(x^2 + (z+1)^2)^{3/2}}, 0, \frac{z-1}{(x^2 + (z-1)^2)^{3/2}} - \frac{z+1}{(x^2 + (z+1)^2)^{3/2}} \right\}$$

Below we plot the charges, equipotential lines, and field lines together:



You should compare this figure with the corresponding one for a (pure) dipole.

Here is a surface plot of $\phi_2(x, z)$ with the equipotential lines superimposed onto the surface:



2.2.8 Quadrupole

It is not hard to extend such computations to arbitrary collections of charges. A combination that is particularly important in the study of nuclear physics, magnets used in particle accelerators, and gravitational waves, is the *quadrupole* which, as its name suggests, consists of 4 poles.

Consider the following arrangement of charges: $+1$ at $\{1, 0, 1\}$ and $\{-1, 0, -1\}$ and -1 at $\{1, 0, -1\}$ and $\{-1, 0, 1\}$:

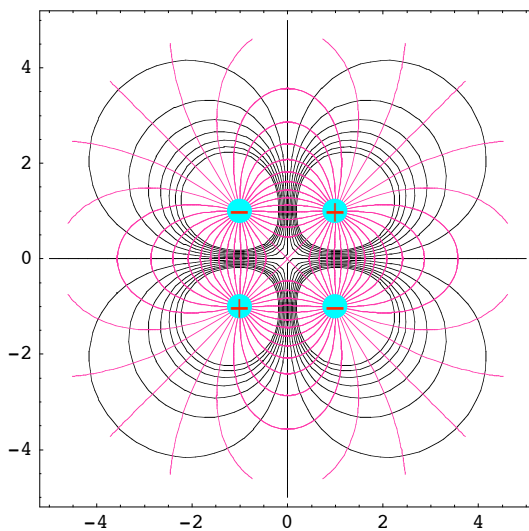
$$\phi_4(\underline{x}, \underline{z}) = \frac{1}{\sqrt{(x+1)^2 + (z-1)^2}} - \frac{1}{\sqrt{(x-1)^2 + (z+1)^2}} + \frac{1}{\sqrt{(x+1)^2 + (z+1)^2}} - \frac{1}{\sqrt{(x-1)^2 + (z-1)^2}}$$

The corresponding electric field is

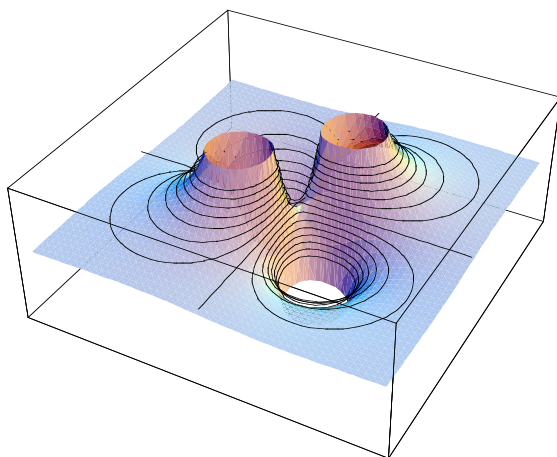
$$\{\mathcal{E}_x, \mathcal{E}_y, \mathcal{E}_z\} = -\nabla\phi_4(x, z)$$

$$\left\{ \begin{aligned} & \frac{x-1}{((x-1)^2 + (z-1)^2)^{3/2}} - \frac{x-1}{((x-1)^2 + (z+1)^2)^{3/2}} - \frac{x+1}{((x+1)^2 + (z-1)^2)^{3/2}} + \frac{x+1}{((x+1)^2 + (z+1)^2)^{3/2}}, 0, \\ & \frac{z-1}{((x-1)^2 + (z-1)^2)^{3/2}} - \frac{z-1}{((x+1)^2 + (z-1)^2)^{3/2}} - \frac{z+1}{((x-1)^2 + (z+1)^2)^{3/2}} + \frac{z+1}{((x+1)^2 + (z+1)^2)^{3/2}} \end{aligned} \right\}$$

Here is a plot of the charges, equipotential lines, and field lines:



Here is a surface plot with the equipotential lines superimposed onto the surface:



Example 2.7 What can you say about the stability of a positive test charge positioned at the origin, $\{0, 0, 0\}$, for the quadrupole potential?

First we need to define *stability*:

Something is *stable* if, after an arbitrary (small) perturbation, the resulting forces acting on it tend to return it to its original position. From the above diagram it is clear that, after a small displacement in the north-east (45°) or south-west (225°) directions, the resulting force would tend to return the test charge to its original position. However, after a small displacement in the north-west (135°) or south-east (315°) directions, the resulting force on the test charge is away from the origin and towards one of the negative charges. Hence a test charge positioned at the origin is not stable.

Note that both the potential and its first (partial) derivatives (*i.e.*, its electric field) at $\{0, 0, 0\}$ are both identically zero:

$$\begin{aligned} &\phi_4(\mathbf{0}, \mathbf{0}) \\ &0 \\ &\{\mathcal{E}_x, \mathcal{E}_y, \mathcal{E}_z\} /. \{x \rightarrow 0, y \rightarrow 0, z \rightarrow 0\} \\ &\{0, 0, 0\} \end{aligned}$$

Using calculus we then know that $\{0, 0, 0\}$ is an *extremum*. In single-variable calculus, if a function has zero derivative then one test to decide whether it is a maximum or a minimum is to compute its second derivative. In higher dimensions there are other (topological) possibilities including *saddle-points*. The generalization of the single-variable test is to compute the eigenvalues of the matrix of second derivatives (*i.e.*, the Hessian):

$$\begin{aligned} &\left(\begin{array}{cc} \frac{\partial^2 \phi_4(x,z)}{\partial x \partial x} & \frac{\partial^2 \phi_4(x,z)}{\partial x \partial z} \\ \frac{\partial^2 \phi_4(x,z)}{\partial z \partial x} & \frac{\partial^2 \phi_4(x,z)}{\partial z \partial z} \end{array} \right) /. \{x \rightarrow 0, y \rightarrow 0, z \rightarrow 0\} \\ &\left(\begin{array}{cc} 0 & \frac{3}{\sqrt{2}} \\ \frac{3}{\sqrt{2}} & 0 \end{array} \right) \end{aligned}$$

If all the eigenvalues are negative (positive) then we have a maximum (minimum). If some of the eigenvalues are positive and some are negative then we have a saddle-point:

$$\begin{aligned} &\mathbf{Eigenvalues}[\%] \\ &\left\{ -\frac{3}{\sqrt{2}}, \frac{3}{\sqrt{2}} \right\} \end{aligned}$$

That we have a saddle-point should be obvious from the surface plot.

2.2.9 Three charges

Three unit positive charges are positioned at the vertices of an equilateral triangle:

$$\mathbf{\Delta} = \left(\begin{array}{ccc} -\frac{1}{2} & 0 & -\frac{1}{2\sqrt{3}} \\ \frac{1}{2} & 0 & -\frac{1}{2\sqrt{3}} \\ 0 & 0 & \frac{1}{\sqrt{3}} \end{array} \right);$$

This problem is dealt with by E Durand in *Electrostatique*, Tome 1, Distributions (Masson, Paris 1964).

Write down the total potential for this configuration of charges;

$$\begin{aligned} &\mathbf{P} = \{x, y, z\}; \\ &\phi_3(x_-, y_-, z_-) = \frac{1}{1} \phi(P, \mathbf{\Delta}[[1]]) + \frac{1}{1} \phi(P, \mathbf{\Delta}[[2]]) + \frac{1}{1} \phi(P, \mathbf{\Delta}[[3]]) \\ &\frac{1}{\sqrt{(x - \frac{1}{2})^2 + y^2 + (z + \frac{1}{2\sqrt{3}})^2}} + \frac{1}{\sqrt{(x + \frac{1}{2})^2 + y^2 + (z + \frac{1}{2\sqrt{3}})^2}} + \frac{1}{\sqrt{x^2 + y^2 + (z - \frac{1}{\sqrt{3}})^2}} \end{aligned}$$

Write down the electric field for this configuration of charges;

$$\{\mathcal{E}_x, \mathcal{E}_y, \mathcal{E}_z\} = -\nabla\phi_3(x, y, z)$$

$$\left\{ \frac{x - \frac{1}{2}}{\left((x - \frac{1}{2})^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}} + \frac{x}{\left(x^2 + y^2 + \left(z - \frac{1}{\sqrt{3}}\right)^2\right)^{3/2}} + \frac{x + \frac{1}{2}}{\left(\left(x + \frac{1}{2}\right)^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}}, \right.$$

$$\frac{y}{\left(x^2 + y^2 + \left(z - \frac{1}{\sqrt{3}}\right)^2\right)^{3/2}} + \frac{y}{\left(\left(x - \frac{1}{2}\right)^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}} + \frac{y}{\left(\left(x + \frac{1}{2}\right)^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}},$$

$$\left. \frac{z - \frac{1}{\sqrt{3}}}{\left(x^2 + y^2 + \left(z - \frac{1}{\sqrt{3}}\right)^2\right)^{3/2}} + \frac{z + \frac{1}{2\sqrt{3}}}{\left(\left(x - \frac{1}{2}\right)^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}} + \frac{z + \frac{1}{2\sqrt{3}}}{\left(\left(x + \frac{1}{2}\right)^2 + y^2 + \left(z + \frac{1}{2\sqrt{3}}\right)^2\right)^{3/2}} \right\}$$

Solve these equations numerically, for example,

```
FindRoot[{E_x, E_y, E_z} == {0, 0, 0}, {{x, 0}, {y, 0}, {z, -0.1}}]
{x -> 0., y -> 0., z -> -0.1643822188}
```

Show that the electric field vanishes at the following 4 points:

$$c = \begin{pmatrix} 0. & 0. & 0. \\ 0. & 0. & -0.1643822188 \\ -0.1423591774 & 0. & 0.08219110939 \\ 0.1423591774 & 0. & 0.08219110939 \end{pmatrix};$$

```
{E_x, E_y, E_z} /. Thread[P -> c[[1]]] // Chop
{0, 0, 0}
```

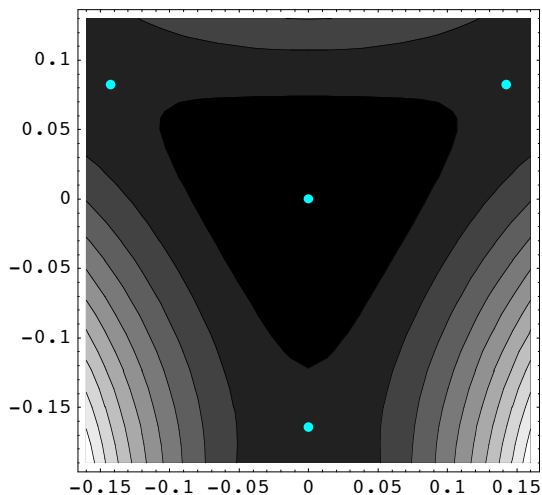
```
{E_x, E_y, E_z} /. Thread[P -> c[[2]]] // Chop
{0, 0, 0}
```

```
{E_x, E_y, E_z} /. Thread[P -> c[[3]]] // Chop
{0, 0, 0}
```

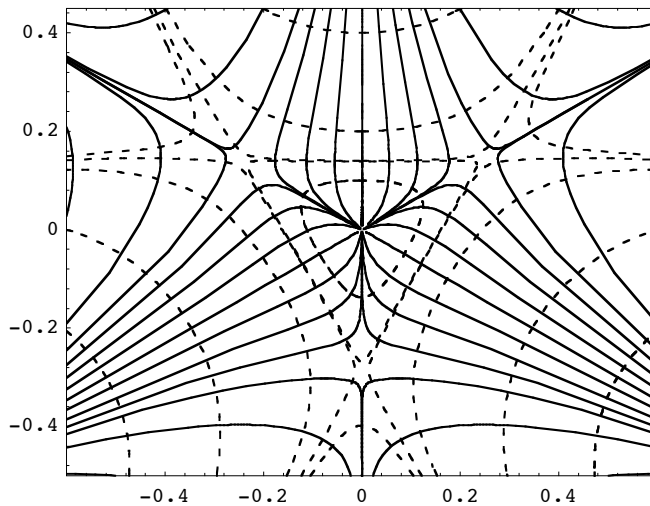
```
{E_x, E_y, E_z} /. Thread[P -> c[[4]]] // Chop
{0, 0, 0}
```

Restricting attention to the x - z plane, here is a plot of the critical points and equipotential contours:

```
ContourPlot[phi_3(x, 0, z), {x, -0.16, 0.16}, {z, -0.19, 0.13},
ContourShading -> True, Epilog -> {PointSize[0.02], Hue[0.5], Point[{0, 0}],
Point[{0, -0.16438}], Point[{-0.142359, 0.08219}], Point[{0.142359, 0.08219}]}];
```



From this plot it looks like $\{0, 0, 0\}$ is a minimum and the other three critical points are saddle-points. Here is a plot of the electric field lines and equipotentials in the x - z plane:



Note that the apparent convergence of flux at the centre is illusory. The flux flow towards the centre diverts out of the plane of the source charges.

What can you say about the stability of a test charge positioned at each of the above 4 points?

We need to compute the Hessian matrix. In 2 dimensions this reads

$$\mathbf{H} = \begin{pmatrix} \frac{\partial^2 \phi_3(x,y,z)}{\partial x \partial x} & \frac{\partial^2 \phi_3(x,y,z)}{\partial x \partial z} \\ \frac{\partial^2 \phi_3(x,y,z)}{\partial z \partial x} & \frac{\partial^2 \phi_3(x,y,z)}{\partial z \partial z} \end{pmatrix};$$

Evaluating the eigenvalues of the Hessian at the first critical point, $\{0, 0, 0\}$:

$H /. Thread[P \to c[1]] // Eigenvalues$
 $\{7.794228634, 7.794228634\}$

it looks like $\{0, 0, 0\}$ is a minimum because both eigenvalues are positive. However, the potential is a function of all three coordinates so we really need to compute the matrix

$$\mathbf{H} = \begin{pmatrix} \frac{\partial^2 \phi_3(x,y,z)}{\partial x \partial x} & \frac{\partial^2 \phi_3(x,y,z)}{\partial x \partial y} & \frac{\partial^2 \phi_3(x,y,z)}{\partial x \partial z} \\ \frac{\partial^2 \phi_3(x,y,z)}{\partial y \partial x} & \frac{\partial^2 \phi_3(x,y,z)}{\partial y \partial y} & \frac{\partial^2 \phi_3(x,y,z)}{\partial y \partial z} \\ \frac{\partial^2 \phi_3(x,y,z)}{\partial z \partial x} & \frac{\partial^2 \phi_3(x,y,z)}{\partial z \partial y} & \frac{\partial^2 \phi_3(x,y,z)}{\partial z \partial z} \end{pmatrix};$$

We now find that all 4 critical points are saddle-points:

$H /. Thread[P \to c[1]] // Eigenvalues$
 $\{-15.58845727, 7.794228634, 7.794228634\}$

$H /. Thread[P \to c[2]] // Eigenvalues$
 $\{24.24375449, -17.07427809, -7.169476398\}$

$H /. Thread[P \to c[3]] // Eigenvalues$
 $\{24.24375449, -17.07427809, -7.169476397\}$

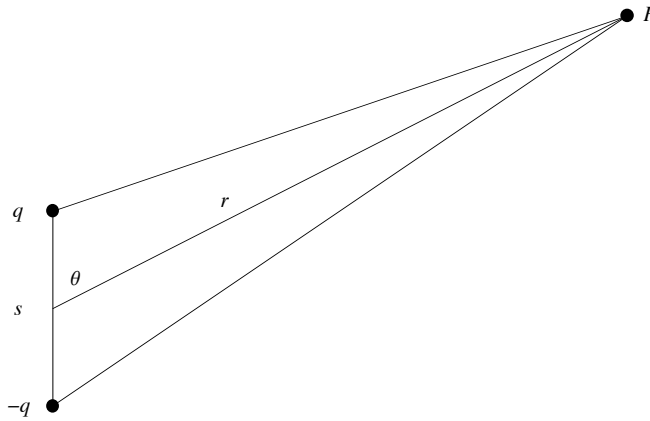
$H /. Thread[P \to c[4]] // Eigenvalues$
 $\{24.24375449, -17.07427809, -7.169476397\}$

That the first critical point (which looked like a minima in two dimensions) is a saddle-point should be obvious from the physical situation: imagine placing a positive test charge at the origin. The force on the test charge after a small displacement out of the plane of the the three fixed charges is *away* from the origin.

The last 3 critical points are themselves vertices of an equilateral triangle. Note that, by symmetry, we should not be suprised then that the eigenvalues of \mathbf{H} evaluated at these critical points are equal.

2.2.10 Multipole Expansion

If you are far away from a localized charge distribution, it "looks" like a point charge, and the potential is approximately $\frac{1}{4\pi\epsilon_0} \frac{Q}{r}$, where Q is the total charge. However, if Q is zero, what is the leading term of the potential for large r ? To answer this question, again consider a physical dipole:



Here the total charge is $Q = q - q = 0$. At the point

$$\mathbf{P} = \{x, y, z\};$$

the potential is

$$\phi_2(\mathbf{x}_-, \mathbf{y}_-, \mathbf{z}_-) = \phi\left(P, \left\{0, 0, \frac{-s}{2}\right\}, -q\right) + \phi\left(P, \left\{0, 0, \frac{s}{2}\right\}, q\right)$$

$$\frac{q}{\sqrt{x^2 + y^2 + \left(z - \frac{s}{2}\right)^2}} - \frac{q}{\sqrt{x^2 + y^2 + \left(\frac{s}{2} + z\right)^2}}$$

In spherical polar coordinates, the potential reads

$$\phi_2(r \cos(\phi) \sin(\theta), r \sin(\phi) \sin(\theta), r \cos(\theta)) // \text{Simplify}$$

$$2q \left(\frac{1}{\sqrt{4r^2 - 4s \cos(\theta)r + s^2}} - \frac{1}{\sqrt{4r^2 + 4s \cos(\theta)r + s^2}} \right)$$

For $r \gg s$ we expand ϕ_2 into a Taylor series in s :

$$\text{Factor} / @ (\% + O[s]^6) // \text{PowerExpand}$$

$$\frac{q \cos(\theta) s}{r^2} + \frac{q \cos(\theta) (5 \cos^2(\theta) - 3) s^3}{8 r^4} + \frac{q \cos(\theta) (63 \cos^4(\theta) - 70 \cos^2(\theta) + 15) s^5}{128 r^6} + O(s^6)$$

The leading term is

$$\phi = \frac{q s \cos(\theta)}{4 \pi \epsilon_0 r^2} = \frac{\mathbf{p} \cdot \hat{\mathbf{r}}}{4 \pi \epsilon_0 r^2}$$

where $p = qs$. This corresponds to a pure dipole potential. Evidently the potential of a dipole goes like $1/r^2$ for large r . Putting together a pair of equal and opposite dipoles makes a *quadrupole*:

$$\begin{aligned} \phi_4(\mathbf{x}_-, \mathbf{y}_-, \mathbf{z}_-) = & \frac{q}{\sqrt{(x - \frac{s}{2})^2 + y^2 + (z - \frac{s}{2})^2}} - \frac{q}{\sqrt{(\frac{s}{2} + x)^2 + y^2 + (z - \frac{s}{2})^2}} - \\ & \frac{q}{\sqrt{(x - \frac{s}{2})^2 + y^2 + (\frac{s}{2} + z)^2}} + \frac{q}{\sqrt{(\frac{s}{2} + x)^2 + y^2 + (\frac{s}{2} + z)^2}} \end{aligned}$$

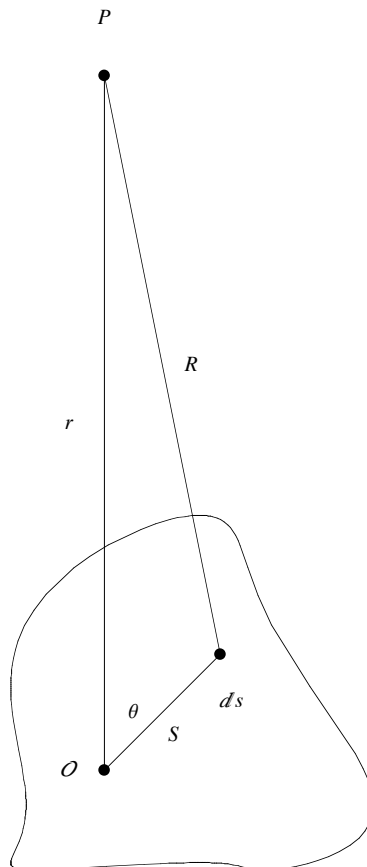
The quadrupole potential goes like $1/r^3$:

$$\begin{aligned} \phi_4(r \cos(\phi) \sin(\theta), r \sin(\phi) \sin(\theta), r \cos(\theta)) + O[s]^4 // \text{Simplify} // \text{PowerExpand} \\ \frac{3q \cos(\theta) \cos(\phi) \sin(\theta) s^2}{r^3} + O(s^4) \end{aligned}$$

The potential of an arbitrary charge distribution confined to a volume V is

$$\phi(\mathbf{P}) = \frac{1}{4\pi\epsilon_0} \int_V \frac{1}{R(s)} \rho(s) ds, \quad (2.5)$$

where R is the distance from ds to \mathbf{P} . With respect to a fixed origin O , we can obtain a systematic expansion for $\phi(\mathbf{P})$ in terms of inverse powers of r . The diagram below defines the variables. Without loss of generality, we have aligned \overline{OP} with the z -axis:



In spherical polar coordinates, we write

$$\mathbf{S} = \{s \cos(\phi) \sin(\theta), s \sin(\phi) \sin(\theta), s \cos(\theta)\};$$

$$\mathbf{P} = \{\mathbf{0}, \mathbf{0}, r\};$$

and find that

$$\mathbf{R} = \|\mathbf{P} - \mathbf{S}\| \text{ // Simplify}$$

$$\sqrt{r^2 - 2s \cos(\theta) r + s^2}$$

For $r > s$ we expand $\frac{1}{R}$ into a series in s :

$$\text{Factor /@ } \left(\frac{1}{R} + O(s^4) \right) \text{ // PowerExpand}$$

$$\frac{1}{r} + \frac{\cos(\theta) s}{r^2} + \frac{(3 \cos^2(\theta) - 1) s^2}{2 r^3} + \frac{\cos(\theta) (5 \cos^2(\theta) - 3) s^3}{2 r^4} + O(s^4)$$

It turns out that the trigonometric terms are *Legendre polynomials*, $P_n(\cos(\theta))$:

$$\text{Factor /@ Table}(P_n(\cos(\theta)), \{n, \mathbf{0}, 3\})$$

$$\left\{ 1, \cos(\theta), \frac{1}{2} (3 \cos^2(\theta) - 1), \frac{1}{2} \cos(\theta) (5 \cos^2(\theta) - 3) \right\}$$

which are generated by the *generating function*:

$$\frac{1}{\sqrt{1 - 2xt + t^2}} = \sum_{n=0}^{\infty} t^n P_n(x), \quad |t| < 1$$

Hence we can write

$$\frac{1}{R(x)} = \frac{1}{r} \sum_{n=0}^{\infty} \left(\frac{s}{r} \right)^n P_n(\cos(\theta)), \quad r > s. \quad (2.6)$$

Combining (2.5) and (2.6) we have

$$\phi(\mathbf{P}) = \frac{1}{4\pi \epsilon_0} \sum_{n=0}^{\infty} \frac{1}{r^{n+1}} \int_V s^n P_n(\cos(\theta)) \rho(s) d\mathbf{s} \equiv$$

$$\frac{1}{4\pi \epsilon_0} \left(\frac{1}{r} \int_V \rho(s) d\mathbf{s} + \frac{1}{r^2} \int_V s \cos(\theta) \rho(s) d\mathbf{s} + \right. \quad (2.7)$$

$$\left. \frac{1}{r^3} \int_V s^2 \frac{1}{2} (3 \cos^2(\theta) - 1) \rho(s) d\mathbf{s} + \dots \right)$$

This is the desired result — the *multipole expansion* of V in powers of $1/r$. The first term ($\sim 1/r$) is the *monopole term*, the second ($\sim 1/r^2$) is the *dipole term*, the third ($\sim 1/r^3$) is the *quadrupole term*, and so on. Although (2.7) is exact is more useful as an approximation scheme. The leading term in the expansion provides the approximate potential at large distances from the charge distribution.

This expansion is not restricted to computing the potential due to a charge distribution: it arises in many fields including atomic and molecular physics (both for bound states of atoms and molecules and in scattering theory), nuclear physics, and gravitational computations.

It is usually easiest to compute (2.7) in spherical polar coordinates. To change coordinates you need to compute the *Jacobian determinant*,

$$d\mathbf{r} \equiv dx dy dz \equiv \begin{vmatrix} \frac{\partial x}{\partial r} & \frac{\partial x}{\partial \theta} & \frac{\partial x}{\partial \phi} \\ \frac{\partial y}{\partial r} & \frac{\partial y}{\partial \theta} & \frac{\partial y}{\partial \phi} \\ \frac{\partial z}{\partial r} & \frac{\partial z}{\partial \theta} & \frac{\partial z}{\partial \phi} \end{vmatrix} dr d\theta d\phi = r^2 \sin(\theta) dr d\theta d\phi, \quad (2.8)$$

In spherical polar coordinates,

$$\mathbf{spc} = r \{\sin(\theta) \cos(\phi), \sin(\theta) \sin(\phi), \cos(\theta)\};$$

the Jacobian matrix reads

$$\begin{pmatrix} \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial r} & \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial \theta} & \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial \phi} \\ \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial r} & \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial \theta} & \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial \phi} \\ \frac{\partial(r \cos(\theta))}{\partial r} & \frac{\partial(r \cos(\theta))}{\partial \theta} & \frac{\partial(r \cos(\theta))}{\partial \phi} \end{pmatrix} \begin{pmatrix} \cos(\phi) \sin(\theta) & r \cos(\theta) \cos(\phi) & -r \sin(\theta) \sin(\phi) \\ \sin(\theta) \sin(\phi) & r \cos(\theta) \sin(\phi) & r \cos(\phi) \sin(\theta) \\ \cos(\theta) & -r \sin(\theta) & 0 \end{pmatrix}$$

Alternatively,

$$D[\mathbf{spc}, \{r, \theta, \phi\}] \begin{pmatrix} \cos(\phi) \sin(\theta) & r \cos(\theta) \cos(\phi) & -r \sin(\theta) \sin(\phi) \\ \sin(\theta) \sin(\phi) & r \cos(\theta) \sin(\phi) & r \cos(\phi) \sin(\theta) \\ \cos(\theta) & -r \sin(\theta) & 0 \end{pmatrix}$$

or

$$\frac{\partial \mathbf{spc}}{\partial \{r, \theta, \phi\}} \begin{pmatrix} \cos(\phi) \sin(\theta) & r \cos(\theta) \cos(\phi) & -r \sin(\theta) \sin(\phi) \\ \sin(\theta) \sin(\phi) & r \cos(\theta) \sin(\phi) & r \cos(\phi) \sin(\theta) \\ \cos(\theta) & -r \sin(\theta) & 0 \end{pmatrix}$$

and the determinant (| |) simplifies to

$$|\%| // \text{Simplify} \\ r^2 \sin(\theta)$$

2.3 Summary

This chapter covers several topics which you should aim to be completely happy with. Here is a summary of these.

Eq. 2.1 was

$$df = \nabla f \cdot d\mathbf{l}$$

and shows how much a function $f(\mathbf{r})$ changes in moving from \mathbf{r} to $\mathbf{r} + d\mathbf{l}$.

If a vector field satisfies

$$\oint_C \mathbf{A} \cdot d\mathbf{l} = 0$$

for *any* circuit C , it is said to be *conservative* and we can write

$$\mathbf{A} = \nabla\psi,$$

where ψ is some scalar function.

In electrostatics the electric field must be conservative and by convention with $\psi = -\phi$ we write

$$\mathbf{E} = -\nabla\phi.$$

Expanding potentials into Taylor series, *e.g.*,

$$\phi(x+h, y+k) = \phi(x, y) + (h \ k) \cdot \begin{pmatrix} \frac{\partial\phi(x,y)}{\partial x} \\ \frac{\partial\phi(x,y)}{\partial y} \end{pmatrix} + \frac{1}{2} (h \ k) \cdot \begin{pmatrix} \frac{\partial^2\phi(x,y)}{\partial x\partial x} & \frac{\partial^2\phi(x,y)}{\partial x\partial y} \\ \frac{\partial^2\phi(x,y)}{\partial y\partial x} & \frac{\partial^2\phi(x,y)}{\partial y\partial y} \end{pmatrix} \begin{pmatrix} h \\ k \end{pmatrix} + \dots,$$

is useful when determining stability and for finding the leading long-range behaviour of a potential.

Chapter 3

Gauss' Law, Gauss' Theorem and Divergence

3.1 Introduction

In this chapter we look at Gauss' Law in a new way. The standard form of Gauss' Law involves integrated quantities *e.g.*, the "flux emergent" from a region is the flux per unit area integrated over the surface. Although this form is very useful in problems with a high degree of symmetry, it only provides a constraint in most other cases without being of much use in finding a functional form for the electric field. In this chapter a *local form* of Gauss' Law is derived that applies at every point.

In proceeding towards the local version of Gauss' Law, a new quantity measuring the production of flux per unit volume is introduced. This scalar quantity is called the *divergence* and can be derived from the field using ∇ , the vector derivative operator of Chapter 2.

3.2 Coulomb's to Gauss' Law

We start with the derivation of Gauss' Law from Coulomb's Law. Gauss' Law contains no new physics beyond Coulomb's Law. Its importance is that it greatly simplifies the problem of finding the electric field in certain cases of simple symmetry. It is also the key to the main work of this chapter which is deriving a differential form of Coulomb's Law. We begin in the standard way by considering the electric flux emerging from a closed surface enclosing a point charge q (Fig. 3.1). The electric flux coming out through an element of area dS equals $E dS \cos(\theta)$ where θ is the angle between the electric field E and a line perpendicular to the area element, dS . It is convenient to think of the surface area element dS as a vector directed along its normal and of magnitude equal to its area, in which the piece of flux can be written as $E \cdot dS$.

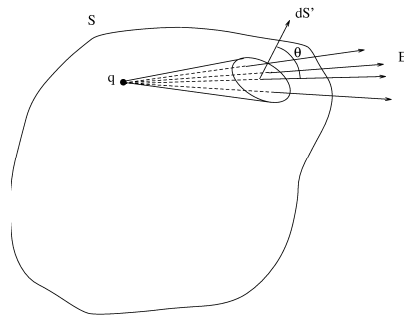


Figure 3.1 The figure shows a surface S which encloses a charge q and a small element of the surface with area dS out of which emerges electric flux.

The total flux emergent from the surface S is then given by

$$\oint_S E \cdot dS,$$

where the circle through the integral sign indicates an integral over a closed surface. In SI units, Coulomb's Law is

$$E = \frac{1}{4\pi\epsilon_0} \frac{q}{r^2} \hat{r},$$

where \hat{r} is a unit vector in the radial direction. Therefore the total flux emergent from the surface is

$$\frac{q}{4\pi\epsilon_0} \oint_S \frac{\hat{r} \cdot d\mathbf{S}}{r^2}.$$

The integrand equals the *projected area* of the element as seen from the point charge (*i.e.*, $\hat{r} \cdot d\mathbf{S}$) divided by its distance squared. This is the definition of the *solid angle* subtended by the element, $d\Omega$. In spherical polar coordinates (see Eq. 2.8):

$$d\mathbf{r} \equiv r^2 d\mathbf{S} = r^2 dr d\Omega, \quad d\mathbf{S} = dS \hat{r} \Rightarrow \frac{\hat{r} \cdot d\mathbf{S}}{r^2} = d\Omega = \sin(\theta) d\theta d\phi.$$

Therefore the total emergent electric flux is

$$\frac{q}{4\pi\epsilon_0} \oint_S d\Omega = \frac{q}{\epsilon_0},$$

where the total solid angle is $\oint_S d\Omega = 4\pi$ steradians.

$$\oint_S d\Omega \equiv \int_0^{2\pi} \int_0^\pi \sin(\theta) d\theta d\phi = 4\pi.$$

Example 3.1: Compute the volume and surface area of a sphere using spherical polar coordinates.

The volume is

$$\frac{\int_0^R \int_0^\pi \int_0^{2\pi} r^2 \sin(\theta) d\phi d\theta dr}{3} = \frac{4\pi R^3}{3}$$

while the surface area of a sphere of radius R is

$$\frac{R^2 \int_0^\pi \int_0^{2\pi} \sin(\theta) d\phi d\theta}{4\pi R^2} = 4\pi R^2$$

Note that the total integral over the solid angle $d\Omega \equiv \sin(\theta) d\phi d\theta$ is

$$\frac{\int_0^\pi \int_0^{2\pi} \sin(\theta) d\phi d\theta}{4\pi} = 4\pi$$

and the SI unit of solid angle is the *steradian*.

Since the electric fields of two charge can be added vectorially, the result can be extended to many charges and we find that the electric flux emergent from a closed surface is equal to the charge enclosed by the surface divided by ϵ_0 . This is Gauss' Law,

$$\oint_S \mathbf{E} \cdot d\mathbf{S} = \frac{q_{\text{enclosed}}}{\epsilon_0}, \quad (3.1)$$

which depends upon the $1/r^2$ nature of Coulomb's Law. Note that gravitational forces obey an equivalent Law:

$$\oint_S \mathbf{g} \cdot d\mathbf{S} = (4\pi G) m_{\text{enclosed}}.$$

3.3 Applying Gauss' Law

Gauss' Law is only directly useful in deriving electric fields in cases of high symmetry. However, it is very quick to apply it in such cases, especially compared to direct application of Coulomb's Law which generally requires the evaluation of some difficult integrals, and the cases for which it is useful are of wide application. There are 3 cases for which Gauss' Law can be applied. They are (1) a plane, (2) a cylinder and (3) a sphere. We now go through each one.

3.3.1 The electric field due to an infinite plane

We wish to know what electric field is produced by an infinite plane charged with $\sigma \text{ C.m}^{-2}$. Although an infinite plane is an impossible idealisation, any surface looks like one if one is close enough to it (*e.g.*, the "Flat Earth"). To solve the problem using Gauss' Law we need to choose a suitable (gaussian) surface enclosing some charge. This should exploit the symmetry of the problem so that electric field is either parallel or perpendicular to the surface. The electric field from an infinite plane must emerge perpendicular to it as there is no preferred direction parallel to the plane. Thus the shape shown in Fig. 3.2 is chosen so that its curved surfaces run parallel to the field and no flux emerges through them. The end caps of area A on the other hand are perpendicular and so a flux $E A$ escapes through each of them.

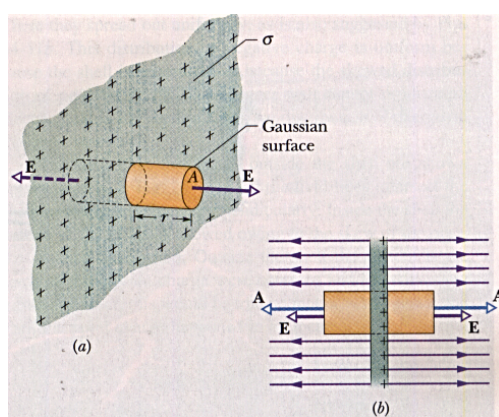


Figure 3.2 The field from an infinite plane emerges at right-angles to it. The surface we consider has arbitrary to faces of the same arbitrary shape and area A which lie parallel to the plane and vertical walls which connect the two faces.

The charge enclosed by this volume is σA and so by Gauss' Law we obtain

$$2EA = \frac{\sigma A}{\epsilon_0},$$

and therefore the magnitude of the electric field from an infinite plane is given by

$$E = \frac{\sigma}{2\epsilon_0}.$$

The field from an infinite plane is equal but opposite on both sides. A more realistic case is the field close to large charged conductor, where "close" implies that it is effectively a plane. This can be treated in exactly the same way except now the field inside the conductor is zero (if it wasn't, current would flow and that would not be electrostatics). Thus all the flux escapes on one side and we get

$$E = \frac{\sigma}{\epsilon_0},$$

for the field close to a charged conductor.

Example 3.2: The electric field beneath a thunder cloud is 1000 V/m. What is the surface charge density of the ground underneath the cloud?

As far as electrostatics are concerned, the Earth is a conductor. Thus $E = \sigma / \epsilon_0$ applies and so $\sigma = 1000 \epsilon_0 = 8.9 \times 10^{-8} \text{ C.m}^{-2}$.

3.3.2 The electric field due to an infinite cylinder

The problem now is to derive the electric field at distance r from the axis of an infinitely long cylinder of radius a charged with $Q \text{ C.m}^{-1}$. The electric field must emerge at right angles to the surface of the cylinder because again there is no preferred direction. Thus the natural gaussian surface is itself a cylinder, but of finite length l as shown in Fig. 3.3.

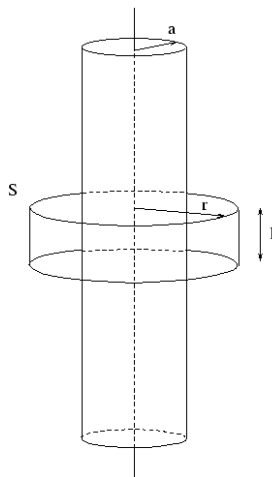


Figure 3.3 The gaussian surface for a long cylinder of radius a is itself a cylinder, but of radius r .

This cylinder is co-axial with the infinite cylinder so that the electric field is uniform over, and perpendicular to, its curved surface. The electric field is parallel to the two ends of the gaussian cylinder and so they do not matter. The surface over which the flux emerges has area $2\pi r l$, while the amount of charge enclosed is $Q l$. Therefore by Gauss' Law we have

$$2\pi r l E = \frac{Q l}{\epsilon_0},$$

and so

$$E = \frac{Q}{2\pi \epsilon_0 r}.$$

Unlike the case of a plane, getting closer to a real cylinder never makes it appear to be an infinite cylinder: end effects do not become infinitesimal. However, there are situations of great practical importance where the above solution is useful. In particular the above field describes the field pattern inside co-axial cables, even in the time-varying case.

3.3.3 The electric field from a charged sphere

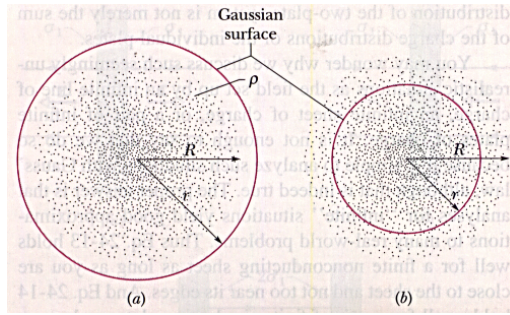
What is the field at a distance r from a sphere with total charge Q (distributed spherically symmetrically)? This case is the easiest. By symmetry the gaussian surface must itself be a sphere centred on the true sphere. The field will come out radially and will therefore be perpendicular to the $4\pi r^2$ area of the gaussian sphere. Thus by Gauss' Law

$$4\pi r^2 E = \frac{Q}{\epsilon_0},$$

and we arrive at the familiar result

$$E = \frac{Q}{4\pi \epsilon_0 r^2}.$$

This is so familiar that it almost seems "obvious" but try deriving it directly from Coulomb's Law and you will see that it is not. This result also applies for an arbitrary spherically symmetric charge distribution where $Q(r)$ is the charge enclosed in a (Gaussian) sphere of radius r



3.3.4 Gauss' Law at a point

To derive a *local version* of Gauss' Law we first need to restate it in mathematical form. We have already seen that the emergent flux can be written $\oint_S \mathbf{E} \cdot d\mathbf{S}$. The charge enclosed can be written as $\int_V \rho d\mathbf{r}$ where ρ is the *charge density* defined throughout the volume V whose bounding surface is $S = \partial V$. Thus Gauss' Law becomes

$$\oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \frac{1}{\epsilon_0} \int_V \rho d\mathbf{r}. \quad (3.2)$$

This is a fundamental equation which you need to remember. Even point charges can be included in this formulation by use of *delta functions*.

We want a version of Gauss' Law that applies at a point. However, one cannot define a volume enclosed or a surface area for a point, and so we consider instead a finite volume that is shrunk to infinitesimal dimensions.

Consider first the charge enclosed

$$Q = \int_V \rho d\mathbf{r},$$

as V becomes smaller. For a continuous charge distribution, there comes a point when V is so small that ρ is essentially constant throughout it and so in the limit $V \rightarrow 0$

$$Q \rightarrow \rho V.$$

We want a finite limit so it makes more sense to divide by V so that we have

$$\lim_{V \rightarrow 0} \frac{1}{V} \int_V \rho d\mathbf{r} = \rho.$$

This leaves us to consider the following limit for the left-hand side of Eq. 3.2, called the *divergence* of the electric field ($\text{div } \mathbf{E}$ for short):

$$\text{div } \mathbf{E} = \lim_{V \rightarrow 0} \frac{1}{V} \oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S}. \quad (3.3)$$

In words this quantity is the amount of electric flux produced per unit volume at a point. It can be defined similarly for any vector field. For instance we will find later that the divergence of the magnetic field is always zero. With this definition Gauss' Law at a point becomes $\text{div } \mathbf{E} = \rho/\epsilon_0$, which says that the amount of electric flux produced per unit volume is proportional to the charge density at every point.

Eq. 3.3 defines the divergence. By considering particular shapes for the volume V , we can obtain expressions for computing the divergence that are suited to particular geometries. Cartesian coordinates are most commonly used, and so let us consider a small cuboid oriented with its sides along the x , y and z axes and centred on the point (x, y, z) :

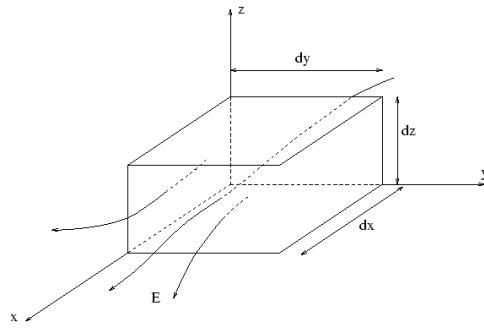


Figure 3.4 A small box with sides parallel to the cartesian axes and drawn to have more electric flux leaving than entering.

Let its sides have lengths Δx , Δy , and Δz . We will now calculate the flux emergent from this cuboid. First consider the amount of flux emerging from the two faces oriented parallel to the y - z plane. Only the x component of the electric field, E_x , contributes to the flux through these faces, and in one face it points in while at the other it points out.

△ A subscripted function such as $G_x(x, y, z)$ denotes the x -component of the vector $\mathbf{G}(x, y, z)$. Partial derivatives are denoted using any of the equivalent standard notations $\partial_x G(x, y, z)$, $\frac{\partial G(x, y, z)}{\partial x}$, or $G^{(1,0,0)}(x, y, z)$. It is mathematically sloppy to denote the partial derivative of a function using a subscript on the function for how would you interpret an expression like $\sum_{i=1}^3 G_i$?

Taking the difference between the x components evaluated in the centre of each face and multiplying by their area these faces contribute

$$\left(E_x \left(x + \frac{\Delta x}{2}, y, z \right) - E_x \left(x - \frac{\Delta x}{2}, y, z \right) \right) \Delta y \Delta z, \quad (3.4)$$

to the flux emergent from the cuboid. The only reason that there is any net contribution to the flux is that the E_x component may change across the cuboid so that the two faces do not cancel. Thus Fig. 3.4 has been drawn to indicate that more flux leaves than enters the box. As Δx becomes small, the expression for E_x can be expanded to first order *e.g.*,

$$E_x \left(x + \frac{\Delta x}{2}, y, z \right) = E_x(x, y, z) + \frac{1}{2} \frac{\partial E_x}{\partial x} \Delta x.$$

The partial derivative applies as the change is in x alone. A similar expression with a negative sign applies for the other face and substituting into Eq. 3.4 we obtain a contribution to the emergent flux of

$$\frac{\partial E_x}{\partial x} \Delta x \Delta y \Delta z.$$

Alternatively, using *Mathematica* we obtain the same result immediately:

$$\left(\mathcal{E}_x \left(x + \frac{\Delta x}{2}, y, z \right) - \mathcal{E}_x \left(x - \frac{\Delta x}{2}, y, z \right) \right) \Delta y \Delta z + O[\Delta x]^2 // \text{Normal} \\ \Delta x \Delta y \Delta z \mathcal{E}_x^{(1,0,0)}(x, y, z)$$

The other four faces give analogous contributions from the y and z components and, recognising the product of lengths $\Delta x \Delta y \Delta z$ as the volume V , we get a total emergent flux from the cuboid of

$$\oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \left(\frac{\partial E_x}{\partial x} + \frac{\partial E_y}{\partial y} + \frac{\partial E_z}{\partial z} \right) V.$$

Therefore the limit of Eq. 3.2, which we called the divergence of \mathbf{E} , becomes

$$\text{div } \mathbf{E} = \lim_{V \rightarrow 0} \frac{1}{V} \oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \frac{\partial E_x}{\partial x} + \frac{\partial E_y}{\partial y} + \frac{\partial E_z}{\partial z}.$$

Using ∇ , the vector derivative operator of Chapter 2, the expression on the right can be written in shorthand form as $\nabla \cdot \mathbf{E}$. We thus arrive at our target, a form of Gauss' Law that applies at a point:

$$\nabla \cdot \mathbf{E} = \frac{\partial E_x}{\partial x} + \frac{\partial E_y}{\partial y} + \frac{\partial E_z}{\partial z} = \frac{\rho}{\epsilon_0}. \quad (3.5)$$

This should be committed to memory as the first of Maxwell's equations. Like all equations, it is best remembered not just as a collection of symbols but from the physical meaning of the various terms. Remembering that $\nabla \cdot \mathbf{E}$ — the divergence of \mathbf{E} — represents the amount of electric flux produced per unit volume, by Gauss' Law it must equal the charge per unit volume, ρ , divided by ϵ_0 .

3.4 Calculating the divergence

△ The divergence of a *vector function* \mathbf{v} is itself a *scalar* $\nabla \cdot \mathbf{v}$. You cannot compute the divergence of a scalar: that is meaningless. In general, the value of the divergence depends on the point at which $\nabla \cdot \mathbf{v}$ is evaluated.

Geometrical interpretation: The name *divergence* should indicate to you that the divergence ($\nabla \cdot \mathbf{v}$) measures how much the vector diverges from the point in question.

If a field \mathbf{v} can be simply expressed in terms of cartesian coordinates, application of Eq. 3.5 is probably the easiest method to compute $\nabla \cdot \mathbf{v}$.

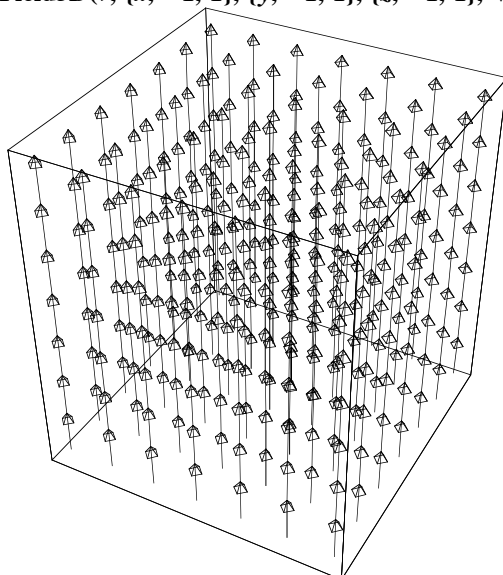
Example 3.3: What is the divergence of the vector function $\mathbf{v} = \hat{z}$?

First, let us visualise this vector field:

$$\mathbf{v} = \{0, 0, 1\};$$

<< Graphics`PlotField3D`

PlotVectorField3D(v, {x, -1, 1}, {y, -1, 1}, {z, -1, 1}, VectorHeads → True);



From the geometrical interpretation we expect this field to have zero divergence at any point $\{x, y, z\}$. Computing the divergence in cartesian coordinates we obtain

$$\frac{\partial v[1]}{\partial x} + \frac{\partial v[2]}{\partial y} + \frac{\partial v[3]}{\partial z}$$

0

Example 3.4: An electric field has the form $E_x = kx$, $E_y = E_z = 0$. What is its divergence and what physical set-up could give such a field?

This is about the simplest possible field other than a constant. We obtain immediately $\nabla \cdot \mathbf{E} = k$:

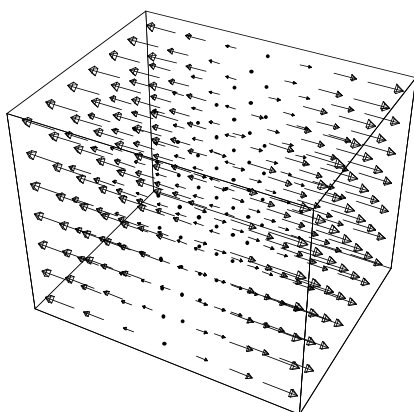
$$\mathcal{E} = \{kx, 0, 0\};$$

$$\frac{\partial \mathcal{E}[1]}{\partial x} + \frac{\partial \mathcal{E}[2]}{\partial y} + \frac{\partial \mathcal{E}[3]}{\partial z}$$

$$k$$

Here is a plot of the vector field:

`PlotVectorField3D(E /. k -> 1, {x, -1, 1}, {y, -1, 1}, {z, -1, 1}, VectorHeads -> True);`



The physical interpretation follows from Gauss' Law. The charge density is proportional to $\nabla \cdot \mathbf{E}$ and so this field comes from a *uniform* charge density and would be the form of field set up inside an infinite slab of uniform charge density perpendicular to the x -axis.

3.4.1 Non-cartesian coordinate systems

It is more difficult if the field is more naturally expressed in a different coordinate system. We had a similar circumstance in Chapter 2 where we showed that the fundamental definition of the gradient could be used to help out. We can do something similar for divergence. Suppose that we have a field of the form

$$\mathbf{E} = E(r) \hat{r}.$$

This is spherically symmetric and in general it is tricky and tedious to apply the Cartesian form of $\nabla \cdot \mathbf{E}$ to such a field. Instead we return to the definition of divergence, Eq. 3.3, but instead of applying it to a cuboid as we did in obtaining $\text{div } \mathbf{E} = \nabla \cdot \mathbf{E}$, we use a shape more suited to the field: we take the small volume V to be a thin spherical shell centred upon the centre of symmetry of the field, with inner and outer radii of r and $r + dr$ respectively.

Given the symmetry of the field and the choice of a shell, whatever the divergence is, it has the same value throughout the shell. Thus the total flux produced by the shell equals the (constant) divergence times the volume of the shell, $V = 4\pi r^2 dr$. The flux produced by the shell, $\Phi = \int_S \mathbf{E} \cdot d\mathbf{S}$, also equals the flux going out through the outer surface, $\Phi_{r+dr} = E(r+dr) A(r+dr) = 4\pi (r+dr)^2 E(r+dr)$ minus that coming in through the inner surface, $\Phi_r = E(r) A(r) = 4\pi r^2 E(r)$, i.e.,

$$\Phi = 4\pi (r+dr)^2 E(r+dr) - 4\pi r^2 E(r) = 4\pi \frac{\partial(r^2 E(r))}{\partial r} dr = \left(\frac{1}{r^2} \frac{\partial(r^2 E(r))}{\partial r} \right) V,$$

where the middle term follows from taking small differences of the expression $r^2 E(r)$ treated as a single function. We can easily verify this result using *Mathematica*:

$$4\pi (r+dr)^2 \mathcal{E}(r+dr) - 4\pi r^2 \mathcal{E}(r) + O(dr)^2 == (4\pi r^2 dr) \left(\frac{1}{r^2} \frac{\partial(r^2 \mathcal{E}(r))}{\partial r} \right) + O(dr)^2$$

True

Hence, from the definition of divergence,

$$\nabla \cdot \mathbf{E} \equiv \lim_{V \rightarrow 0} \frac{1}{V} \oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S},$$

we obtain

$$\nabla \cdot \mathbf{E} = \frac{1}{r^2} \frac{\partial(r^2 E_r)}{\partial r}.$$

Proceeding in this manner one can obtain more general expressions for the divergences of fields expressed in spherical polar and other coordinates. It turns out that there are much more direct methods for computing the divergence in any coordinate system — see Appendix A for a summary of vector operators in orthogonal coordinate systems. For example, the general expression for the divergence in spherical polar coordinates is (see Eq. A.14)

$$\nabla \cdot \mathbf{E} = \frac{1}{r^2} \frac{\partial}{\partial r} (r^2 E_r) + \frac{1}{r \sin(\theta)} \frac{\partial}{\partial \theta} (\sin(\theta) E_\theta) + \frac{1}{r \sin(\theta)} \frac{\partial E_\phi}{\partial \phi}, \quad (3.6)$$

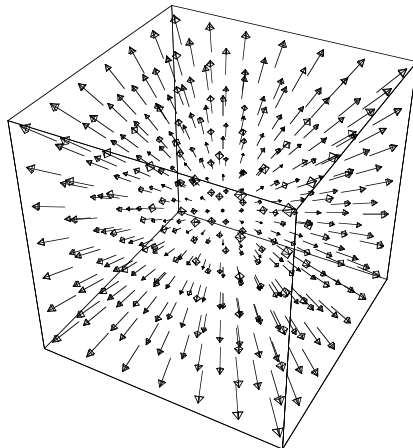
where $\mathbf{E} \equiv E_r \hat{\mathbf{r}} + E_\theta \hat{\boldsymbol{\theta}} + E_\phi \hat{\boldsymbol{\phi}}$.

Example 3.5: What is the divergence of the vector function $\mathbf{v} = \mathbf{r} = r \hat{\mathbf{r}}$?

First, let us visualise this vector field using cartesian coordinates:

$$\mathbf{v} = \{x, y, z\};$$

`PlotVectorField3D(v, {x, -1, 1}, {y, -1, 1}, {z, -1, 1}, VectorHeads → True);`



From the geometrical interpretation we expect this field to have large (positive) divergence at any point $\{x, y, z\}$. Computing the divergence in cartesian coordinates we obtain

$$\frac{\partial v[1]}{\partial x} + \frac{\partial v[2]}{\partial y} + \frac{\partial v[3]}{\partial z}$$

3

Since $\mathbf{v} \equiv v_r \hat{\mathbf{r}} + v_\theta \hat{\boldsymbol{\theta}} + v_\phi \hat{\boldsymbol{\phi}} = r \hat{\mathbf{r}}$, using the divergence in spherical polar coordinates (Eq. 3.6) we find that

$$\frac{1}{r^2} \frac{\partial(r^2 r)}{\partial r}$$

3

which is identical to the result obtained using cartesian coordinates. For this particular vector field, the divergence does not depend on the point at which it is computed.

Example 3.6: What is the divergence of $\mathbf{v} = \frac{1}{r^2} \hat{\mathbf{r}} \equiv \frac{1}{r^3} \mathbf{r}$?

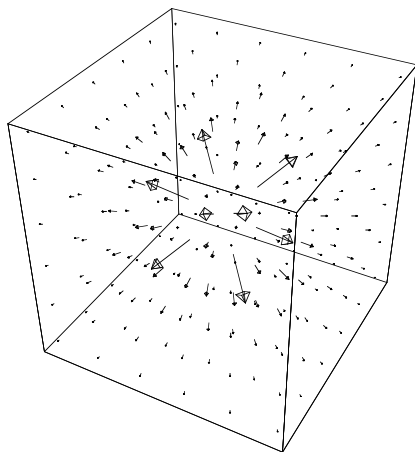
With

$$\mathbf{v} = \frac{\{x, y, z\}}{(x^2 + y^2 + z^2)^{3/2}}$$

$$\left\{ \frac{x}{(x^2 + y^2 + z^2)^{3/2}}, \frac{y}{(x^2 + y^2 + z^2)^{3/2}}, \frac{z}{(x^2 + y^2 + z^2)^{3/2}} \right\}$$

we visualize the field:

PlotVectorField3D(v, {x, -1, 1}, {y, -1, 1}, {z, -1, 1}, VectorHeads → True, PlotPoints → 6);



From this plot we would expect this field to have non-zero divergence. However, computing the divergence in cartesian coordinates,

$$\frac{\partial v[1]}{\partial x} + \frac{\partial v[2]}{\partial y} + \frac{\partial v[3]}{\partial z} \text{ // Together}$$

$$0$$

or spherical polar coordinates,

$$\frac{1}{r^2} \partial_r \left(r^2 \frac{1}{r^2} \right)$$

$$0$$

we find that the divergence is identically zero! What is going on here? We will return to this example shortly.

3.5 Poisson's equation, Laplace's equation, and Gauss' Theorem

3.5.1 Poisson's equation

We showed in Chapter 2 that conservation of energy means that an electrostatic field can be expressed in terms of a potential as in $\mathbf{E} = -\nabla\phi$. Substituting this into Eq. 3.5 we obtain

$$\nabla \cdot \nabla\phi \equiv \Delta\phi = \nabla^2\phi = -\frac{\rho}{\epsilon_0}, \quad (3.7)$$

which is known as *Poisson's equation*. Here ∇^2 is the *Laplacian operator* which, written in Cartesian form, is

$$\nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2}.$$

The Laplacian operator also arises in quantum mechanics. There are other forms for ∇^2 in different coordinate systems. *E.g.*, the Laplacian operator in spherical polar coordinates reads (Eq. A.15)

$$\nabla^2 = \frac{1}{r^2} \frac{\partial}{\partial r} \left(r^2 \frac{\partial}{\partial r} \right) + \frac{1}{r^2} \left(\frac{1}{\sin(\theta)} \frac{\partial}{\partial \theta} \left(\sin(\theta) \frac{\partial}{\partial \theta} \right) + \frac{1}{\sin^2(\theta)} \frac{\partial^2}{\partial \phi^2} \right). \quad (3.8)$$

Poisson's equation (Eq. 3.7) can be used to find the charge distribution given a form for the potential.

Example 3.7: What charge distribution is needed to give a potential of the form $\phi = k r^2$ where r is the distance from a point?

Apply the ∇^2 operator to $r^2 = x^2 + y^2 + z^2$. Thus

$$\begin{aligned} \frac{\partial}{\partial x} r^2 &= 2x, \\ \frac{\partial}{\partial x} 2x &= 2, \end{aligned}$$

$$\text{so } \nabla^2 (k r^2) = 6k.$$

Alternatively, using the Laplacian operator in spherical polar coordinates, we get the same result:

$$\nabla^2 (k r^2) = k \frac{1}{r^2} \frac{\partial}{\partial r} \left(r^2 \frac{\partial(r^2)}{\partial r} \right) = k \frac{1}{r^2} \frac{\partial}{\partial r} (2r^3) = 6k.$$

Therefore Eq. 3.7 gives $\rho = -6k \epsilon_0$. Thus a constant charge density gives a potential proportional to r^2 and this is the form of potential inside a uniformly charged sphere for example.

3.5.2 Laplace's equation

In regions with no charge density, Poisson's equation reduces to

$$\nabla^2 \phi = 0, \quad (3.9)$$

which is known as *Laplace's equation*. Solutions of this equation with boundary conditions are important in the design of the focussing fields of TV tubes for instance.

Example 3.8: Verify that the $1/r$ Coulomb potential satisfies Laplace's equation.

Using brute force by applying the ∇^2 operator in cartesian form, *i.e.*,

$$\nabla^2 = \frac{\partial^2}{\partial x^2} + \frac{\partial^2}{\partial y^2} + \frac{\partial^2}{\partial z^2},$$

we have to calculate derivatives such as

$$\frac{\partial^2 (1/r)}{\partial x^2}.$$

Since $r^2 = x^2 + y^2 + z^2$ we have

$$\frac{\partial(1/r)}{\partial x} = -\frac{1}{r^2} \frac{\partial r}{\partial x} = -\frac{x}{r^3}.$$

Computing the second derivative

$$\frac{\partial(-x/r^3)}{\partial x} = -\frac{1}{r^3} + \frac{3x}{r^4} \frac{\partial r}{\partial x} = -\frac{1}{r^3} + \frac{3x^2}{r^5}.$$

Similar expressions apply to the other components and we find

$$\nabla^2 \left(\frac{1}{r} \right) = -\frac{3}{r^3} + \frac{3(x^2 + y^2 + z^2)}{r^5} = 0,$$

since $x^2 + y^2 + z^2 = r^2$.

Alternatively, using the Laplacian operator in spherical polar coordinates (Eq. A.15),

$$\nabla^2\left(\frac{1}{r}\right) = \frac{1}{r^2} \frac{\partial}{\partial r} \left(r^2 \frac{\partial(1/r)}{\partial r} \right) = \frac{1}{r^2} \frac{\partial}{\partial r} (-1) = 0.$$

3.5.3 Gauss' Theorem

The interpretation of divergence as the flux produced per unit volume suggests that the following integral

$$\int_V \nabla \cdot \mathbf{A} \, d\mathbf{r},$$

must represent the emergent flux from a volume V for *any* vector field \mathbf{A} . We have seen that the emergent flux can also be written as $\oint_{S=\partial V} \mathbf{A} \cdot d\mathbf{S}$ and so therefore we expect

$$\oint_{S=\partial V} \mathbf{A} \cdot d\mathbf{S} = \int_V \nabla \cdot \mathbf{A} \, d\mathbf{r}. \quad (3.10)$$

This result is known as *Gauss' Theorem* (sometimes it is called the *divergence theorem*). It is important to distinguish between *Gauss' Theorem*, which has only mathematical content, and applies to *any* physical vector field, as opposed to *Gauss' Law* which is founded in experiment and is just another way of expressing Coulomb's Law.

The importance of Gauss' Theorem is that it provides a way to transform between the surface and volume integrals frequently encountered in physics. Thus if we go back to the integral version of Gauss' Law (Eq. 3.1)

$$\oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \frac{1}{\epsilon_0} \int_V \rho \, d\mathbf{r},$$

and apply Gauss' Theorem, we can immediately deduce that

$$\oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \int_V \nabla \cdot \mathbf{E} \, d\mathbf{r} = \frac{1}{\epsilon_0} \int_V \rho \, d\mathbf{r},$$

Since this applies for *any* volume V , we must have

$$\nabla \cdot \mathbf{E} = \frac{\rho}{\epsilon_0},$$

which, as before, is Gauss' Law in differential form (but it no longer depends upon the assumption of a volume of particular shape as that is accounted for in the proof of Gauss' Theorem).

3.5.4 A paradox — the Dirac delta function

Example 3.9: Compute $\nabla \cdot \mathbf{E}$ for a point charge and compare with ρ/ϵ_0 .

The potential of a point charge, Q , is

$$\phi(r) = \frac{1}{4\pi\epsilon_0} \frac{Q}{r}.$$

Hence the electric field \mathbf{E} is

$$\mathbf{E}(r) = -\nabla\phi(r) = \frac{-\partial\phi(r)}{\partial r} \hat{\mathbf{r}} = \frac{1}{4\pi\epsilon_0} \frac{Q}{r^2} \hat{\mathbf{r}}.$$

In Example 3.6 we saw that

$$\mathbf{v} = \frac{1}{r^2} \hat{\mathbf{r}} \Rightarrow \nabla \cdot \mathbf{v} = 0 \Rightarrow \nabla \cdot \mathbf{E} = 0.$$

However, after visualizing the fields we were puzzled to find that the divergence was identically zero.

Alternatively, in Example 3.8 we found that $\nabla^2\left(\frac{1}{r}\right) = 0$. Hence

$$\nabla \cdot \mathbf{E} \equiv -\nabla^2 \phi = -\frac{Q}{4\pi\epsilon_0} \nabla^2\left(\frac{1}{r}\right) = 0.$$

Gauss' Law in differential form says that

$$\nabla \cdot \mathbf{E} = \frac{\rho}{\epsilon_0},$$

and, for a point charge, the charge density is zero everywhere except at $r = 0$ where it is infinite!

If we apply the divergence theorem to the electric field \mathbf{E} we find that

$$\int_V \nabla \cdot \mathbf{E} \, d\mathbf{r} = \oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} = \frac{Q}{4\pi\epsilon_0} \oint_{S=\partial V} \frac{1}{r^2} r^2 \, d\Omega = \frac{Q}{\epsilon_0}.$$

However, above we have shown that $\nabla \cdot \mathbf{E} = 0$. What is the resolution to this paradox? *Hint:* you should be suspicious of "point" charges. Taking Coulomb's Law at face value, the potential and field of a "point charge" at the origin are infinite. Note that, although there is no such thing as a "point charge", the electron is *effectively* a point charge with physical radius $\lesssim 10^{-17} \text{ m}$.

The problem is the point $r = 0$, where \mathbf{E} blows up. A more careful analysis shows that $\nabla \cdot \mathbf{E} = 0$ everywhere *except* at the origin. We seem to require a function with the bizarre property that $\nabla \cdot \mathbf{E} = 0$ everywhere except at a single point, yet

$$\int_V \nabla \cdot \mathbf{E} \, d\mathbf{r} = \frac{Q}{\epsilon_0}.$$

No function can possibly behave this way. What we have stumbled onto is the *Dirac delta function* which is not really a function at all. The Dirac delta "function" was originally defined by Dirac as

$$\delta(x) = \begin{cases} 0, & x \neq 0 \\ \infty, & x = 0 \end{cases}, \text{ and } \int_{-\infty}^{\infty} \delta(x) \, dx = 1. \quad (3.11)$$

However this definition does not make sense mathematically. In addition, $\delta(x)$ has the interesting property that

$$\int_{-\infty}^{\infty} f(x-a) \delta(x) \, dx = f(a). \quad (3.12)$$

It turns out that Dirac had a good idea though. See Appendix B for more on $\delta(x)$.

Example 3.10: Compute \mathbf{E} , ϕ , and $\nabla \cdot \mathbf{E}$ for the following (spherically symmetric) charge distribution:

$$\rho(\mathbf{r}) = \begin{cases} \frac{Q}{\frac{4}{3}\pi R^3}, & r < R \\ 0, & r \geq R \end{cases}.$$

Using spherical polar coordinates, the charge enclosed in a sphere of radius $r < R$ is

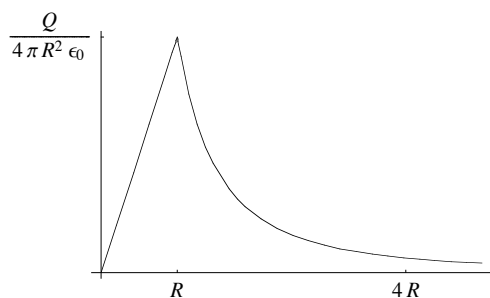
$$q(r) = \int_V \rho(\mathbf{r}) \, d\mathbf{r} = 4\pi \int_0^r \rho(r) r^2 \, dr = 4\pi \frac{Q}{\frac{4}{3}\pi R^3} \int_0^r r^2 \, dr = 4\pi \frac{Q}{\frac{4}{3}\pi R^3} \frac{r^3}{3} = Q \frac{r^3}{R^3}.$$

For $r = R$, we find that $q(R) = Q$, the *total charge*.

Using Gauss' Law (also see Section 3.3.3), the electric field is

$$\mathbf{E}(r) = \begin{cases} \frac{1}{4\pi\epsilon_0} \frac{Qr}{R^3} \hat{\mathbf{r}}, & r < R \\ \frac{1}{4\pi\epsilon_0} \frac{Q}{r^2} \hat{\mathbf{r}}, & r \geq R \end{cases}.$$

Here is a plot of the radial component of $\mathbf{E}(r)$, i.e., E_r :



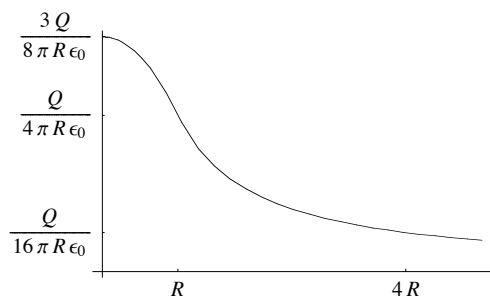
The Gradient Theorem says that

$$\phi(B) - \phi(A) = \int_A^B d\phi = - \int_A^B \mathbf{E} \cdot d\mathbf{l} \Rightarrow \phi(r) = - \int_{\infty}^r E_r \cdot dr,$$

and leads to

$$\phi(r) = \begin{cases} \frac{1}{4\pi\epsilon_0} \frac{Q}{R^3} \left(\frac{3}{2} R^2 - \frac{1}{2} r^2 \right) & , r < R \\ \frac{1}{4\pi\epsilon_0} \frac{Q}{r} & , r \geq R \end{cases}.$$

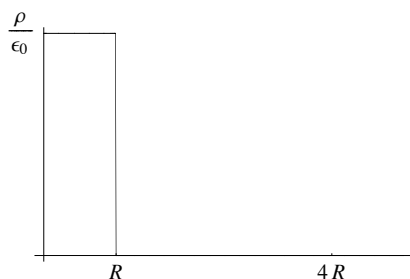
Here is a plot of the potential, $\phi(r)$:



The divergence of the electric field is (see Eq. 3.6)

$$\nabla \cdot \mathbf{E}(r) = \frac{1}{r^2} \frac{\partial}{\partial r} (r^2 E_r) = \begin{cases} \frac{Q}{4\pi\epsilon_0} \frac{3}{R^3} = \frac{1}{\epsilon_0} \left(\frac{Q}{\frac{4}{3}\pi R^3} \right) = \frac{\rho}{\epsilon_0} & , r < R \\ 0 & , r \geq R \end{cases}.$$

Here is a plot of $\nabla \cdot \mathbf{E}(r)$:



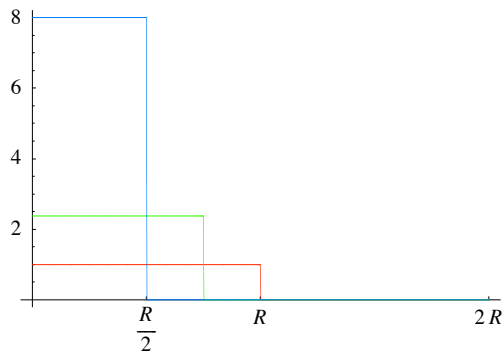
This result is to be expected from Gauss' Law. Note that $\nabla \cdot \mathbf{E}$ is *discontinuous* (because ρ itself is discontinuous).

If we apply the divergence theorem to the electric field \mathbf{E} we find that

$$\begin{aligned} \int_V \nabla \cdot \mathbf{E} d\mathbf{r} &= \int_V \frac{\rho(\mathbf{r})}{\epsilon_0} d\mathbf{r} = \frac{q(r)}{\epsilon_0} \equiv \oint_{S=\partial V} \mathbf{E} \cdot d\mathbf{S} \\ &= \\ \frac{1}{4\pi\epsilon_0} \begin{cases} \oint \frac{Qr}{R^3} r^2 d\Omega & r < R \\ \oint \frac{Q}{r^2} r^2 d\Omega & r \geq R \end{cases} &= \frac{Q}{\epsilon_0} \begin{cases} \frac{r^3}{R^3} & r < R \\ 1 & r \geq R \end{cases} \equiv \frac{q(r)}{\epsilon_0}, \end{aligned}$$

which all checks out. There is no paradox here.

Consider now what happens to $\rho(r) = \frac{Q}{\frac{4}{3}\pi R^3}$ for this spherically symmetric charge distribution in the "point charge" limit, *i.e.*, as $R \rightarrow 0$:



The boxes get narrower and taller — in such a way that the *volume integral* (note the $r^2 dr$ factor) is a constant equal to the total charge. If you read through Appendix B you should be able to show that, in the limit, $\rho(\mathbf{r}) = Q\delta(\mathbf{r})$ and hence $\nabla \cdot \mathbf{E} = \frac{\rho(\mathbf{r})}{\epsilon_0} = \frac{Q}{\epsilon_0}\delta(\mathbf{r})$. This simple result is a key ingredient to a concise formulation much of electromagnetism.

3.5.5 Continuity Equations

As another example of Gauss' Theorem we will look at how the conservation of charge can be expressed as a differential equation. Our result will be used later in deducing the existence of an extra term in one of Maxwell's equations, a term that allows the propagation of electromagnetic waves. Since there are conserved quantities in many branches of physics, equations of very similar form crop up over and over again and are known as *continuity equations*.

Consider the charge flowing out of a volume V . The total rate of charge leaving V is given by the integral of the current density \mathbf{J} over the surface of the volume. \mathbf{J} is a vector with dimensions $A.m^{-2}$ directed along the local current flow at every point and with magnitude equal to the current density. Thus if the total charge in V is Q , the conservation of charge tells us that the charge flowing out of V must be balanced by a decrease in Q , that is

$$\frac{\partial Q}{\partial t} = -\oint_S \mathbf{J} \cdot d\mathbf{S},$$

where the partial derivative indicates that the volume is fixed in position. Since the total charge $Q = \int_V \rho d\mathbf{r}$ we find

$$\frac{\partial}{\partial t} \int_V \rho d\mathbf{r} = -\oint_S \mathbf{J} \cdot d\mathbf{S}.$$

Now apply Gauss' Theorem to transform the surface integral into a volume integral and we find

$$\int_V \frac{\partial \rho}{\partial t} d\mathbf{r} = -\int_V \nabla \cdot \mathbf{J} d\mathbf{r}.$$

As before, since this applies for *any* volume V we obtain our final result, the continuity equation for electric charge

$$\frac{\partial \rho}{\partial t} + \nabla \cdot \mathbf{J} = 0. \quad (3.13)$$

This equation expresses the conservation of electric charge. It says that at every point the electric current produced per unit volume ($\nabla \cdot \mathbf{J}$) must be balanced by a decrease in the charge density.

3.5.6 Other examples of the continuity equation

In fluid flow, the equivalent of current density is the mass flow rate per unit area, $\rho_m \mathbf{v}$, where \mathbf{v} is the velocity vector field. The mass density ρ_m takes the place of the charge density, and the continuity equation becomes

$$\frac{\partial \rho_m}{\partial t} + \nabla \cdot (\rho_m \mathbf{v}) = 0.$$

In incompressible flow (a good approximation at low speeds), ρ_m is constant and we have $\nabla \cdot \mathbf{v} = 0$, an important equation in fluid dynamics.

In the conduction of heat in a uniform solid the "density" of heat is CT where C is the heat capacity per unit volume and T is the temperature. The equation of continuity is then

$$C \frac{\partial T}{\partial t} + \nabla \cdot \mathbf{Q} = 0.$$

where \mathbf{Q} is the heat flux. We obtained an expression for \mathbf{Q} in terms of the temperature gradient in Eq. 2.2 and substituting this we obtain

$$\nabla^2 T = \frac{C}{\kappa} \frac{\partial T}{\partial t},$$

a fundamental equation in the theory of heat conduction, also known as the *diffusion equation*.

3.6 Summary

In this chapter we expressed Gauss' Law in integral form as

$$\oint_S \mathbf{E} \cdot d\mathbf{S} = \frac{1}{\epsilon_0} \int_V \rho \, d\mathbf{r},$$

We then progressed from that to considering the limit of infinitesimal volumes, defining a scalar quantity called the divergence $\text{div } \mathbf{E}$ by

$$\text{div } \mathbf{E} = \lim_{V \rightarrow 0} \frac{1}{V} \oint_S \mathbf{E} \cdot d\mathbf{S}.$$

A form convenient for cartesian coordinates was developed by considering a volume in the shape of a cuboid. We obtained

$$\text{div } \mathbf{E} = \frac{\partial E_x}{\partial x} + \frac{\partial E_y}{\partial y} + \frac{\partial E_z}{\partial z} = \nabla \cdot \mathbf{E}.$$

Gauss' Law at a point was finally derived:

$$\nabla \cdot \mathbf{E} = \frac{\rho}{\epsilon_0}.$$

We then returned to the Gauss' Theorem, a mathematical theorem that applies to any continuous vector field and allows one to transform between surface and volume integrals

$$\oint_S \mathbf{A} \cdot d\mathbf{S} = \int_V \nabla \cdot \mathbf{A} \, d\mathbf{r}.$$

Gauss' theorem was applied to derive the continuity equation which expresses the conservation of charge in differential form:

$$\frac{\partial \rho}{\partial t} + \nabla \cdot \mathbf{J} = 0,$$

where \mathbf{J} is the current density. Finally the examples illustrated how to cope with non-cartesian coordinate systems.

Chapter 4

Faraday's Law, Stokes' Theorem and curl

4.1 Introduction

This chapter repeats the pattern of Chapter 3. We start from an experimentally derived physical law, in this case Faraday's Law of induction, and derive a differential version of it that applies at a point. In so doing, we introduce the third and final of the vector derivatives, a vector called the *curl* of a field. We then follow up with Stokes' Theorem which, in the same way that Gauss' Theorem is used to transform between volume and surface integrals, can be used to transform between surface and line integrals.

4.2 Faraday's Law of Induction

Faraday observed that changing the magnetic flux through a loop of wire whether by moving the wire or the source of the magnetic field caused a voltage to be developed around the loop. The voltage produced is proportional to the rate of change of the magnetic flux through the loop. The direction of the voltage produced is such that if a current flows it "tries" to keep the field constant. This is Lenz's Law and leads to minus signs in the equations for induction. As an aside, super-conducting loops are able to keep the flux *precisely* constant over long periods of time.

Faraday's work is the classic example of basic physics with applications of immense importance (dynamoes, transformers, etc), unrecognised at the time of its discovery.

We start by writing a mathematical version of Faraday's' Law. The voltage, V , or EMF (electro-motive force) around a circuit, C , is simply the line integral of the electric field:

$$V = \oint_C \mathbf{E} \cdot d\mathbf{l}.$$

We met this earlier in Chapter 2 where we said that this quantity had to be zero for energy conservation. However, that was in electrostatics, and does not apply when work is being done to change the fields. A corollary is that the electrostatic relation $\mathbf{E} = -\nabla\phi$ no longer applies in the time-varying case. The flux connecting the circuit C is the integral of the magnetic flux density \mathbf{B} over any surface, S , bounded by $C = \partial S$ and can be written

$$\int_S \mathbf{B} \cdot d\mathbf{S}.$$

Therefore Faraday's Law in integral form is

$$\oint_{C=\partial S} \mathbf{E} \cdot d\mathbf{l} = -\frac{d}{dt} \int_S \mathbf{B} \cdot d\mathbf{S}. \quad (4.1)$$

For the sign to make sense, the direction in which the circuit is travelled has to be defined. Fig. 4.1 shows the convention based upon the right-hand rule. If one grasps the circuit with the right-hand so that the fingers point along the direction of \mathbf{B} , then the thumb points along the direction in which C is traversed.

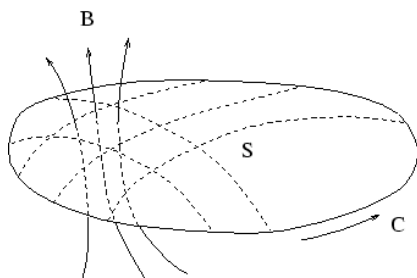


Figure 4.1 Magnetic flux threads a circuit C which is covered by a surface S that has C as its boundary. The arrow on C indicates in which direction the line integral is taken for \mathbf{B} pointing in the direction shown.

Example 4.1: What is the electric field inside a long solenoid of n turns/unit length when the current I flowing through the coils changes?

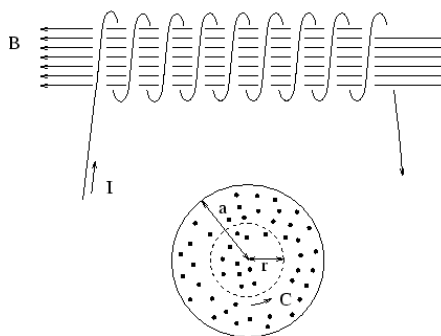


Figure 4.2 The figure shows side and end-on views of a solenoid carrying current I . The end-on view looks into the magnetic field (represented by dots). To calculate the electric field induced by changing I , a circuit is taken to be a circle of radius r enclosing the field.

We will assume the result from first year that the magnetic field inside the solenoid is given by $B = \mu_0 n I$. By symmetry \mathbf{E} must run in circles around the axis of the solenoid and so we take such a circle as our circuit. Since \mathbf{E} runs parallel to the circuit at all points, the line integral reduces to

$$\oint_C \mathbf{E} \cdot d\mathbf{l} = 2\pi r E.$$

For circuits inside the radius a of the solenoid ($r < a$), the flux linking the circuit is

$$\Phi_B = \int_S \mathbf{B} \cdot d\mathbf{S} = \pi r^2 B = \pi r^2 \mu_0 n I.$$

Therefore applying Faraday's Law we have

$$2\pi r E = -\pi r^2 \mu_0 n \frac{dI}{dt},$$

or

$$E = -\frac{r}{2} \mu_0 n \frac{dI}{dt} \quad (r < a).$$

There is no magnetic flux outside the solenoid so the flux linking the circuit stays fixed at $\pi a^2 B$ for $r > a$ and therefore

$$E = -\frac{a^2}{2r} \mu_0 n \frac{dI}{dt} \quad (r > a).$$

The existence of an electric field outside the coil allows signals flowing through the coil to be picked up with a loop of wire enclosing the coil.

4.3 Curl and Stokes' Theorem

As for Gauss' Law, we would like a version of Faraday's Law that applies at a point. We will start by considering the line integral around a loop as the loop is shrunk to infinitesimal size.

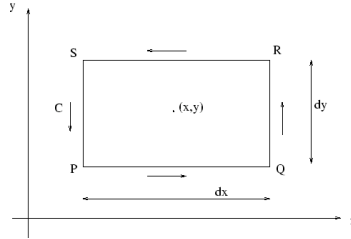


Figure 4.3 The figure shows a small loop used to obtain an expression for the line integral in the limit of infinitesimal size. The circuit is traversed in a direction appropriate for the right-hand rule and a z -axis which points out of the page.

To be specific we consider the loop illustrated in Fig. 4.3. This is a rectangle in the x - y plane with sides parallel to the x and y -axes. Note that by contrast with the derivation of divergence, the loop's orientation *is* significant. This will be reflected in the quantity called *curl* which we will introduce which turns out to be a vector rather than a scalar like divergence. The circuit direction indicated in Fig. 4.3 follows the right-hand rule for a right-handed coordinate set of axes in which $\hat{z} = \hat{x} \wedge \hat{y}$.

The line integral around the circuit has four separate parts corresponding to the line segments \overline{PQ} , \overline{QR} , \overline{RS} and \overline{SP} . The contribution from \overline{PQ} is due entirely to the x component of E , which we evaluate at the mid-point of the segment as $E_x(x, y - dy/2)$. Multiplying this by the length of the segment and adding in the other three similar terms we have

$$\oint_C \mathbf{E} \cdot d\mathbf{l} = E_x(x, y - dy/2) dx + E_y(x + dx/2, y) dy - E_x(x, y + dy/2) dx - E_y(x - dx/2, y) dy,$$

with the minus signs appearing when we travel against the direction of the coordinate axes. This expression can be grouped into two pairs of differences:

$$\oint_C \mathbf{E} \cdot d\mathbf{l} = (E_y(x + dx/2, y) - E_y(x - dx/2, y)) dy - (E_x(x, y + dy/2) - E_x(x, y - dy/2)) dx,$$

which, when expanded to first order as we did when deriving $\text{div } \mathbf{E} \equiv \nabla \cdot \mathbf{E}$, yields:

$$\left(\mathcal{E}_y \left(x + \frac{dx}{2}, y \right) - \mathcal{E}_y \left(x - \frac{dx}{2}, y \right) \right) dy + O(dx)^3 \\ dy \mathcal{E}_y^{(1,0)}(x, y) dx + O((dx)^3)$$

and similarly

$$\left(\mathcal{E}_x \left(x, y + \frac{dy}{2} \right) - \mathcal{E}_x \left(x, y - \frac{dy}{2} \right) \right) dx + O(dy)^3 \\ dx \mathcal{E}_x^{(0,1)}(x, y) dy + O((dy)^3)$$

Collecting terms together, we can write

$$\oint_C \mathbf{E} \cdot d\mathbf{l} = \left(\frac{\partial E_y}{\partial x} - \frac{\partial E_x}{\partial y} \right) dx dy = (\nabla \wedge \mathbf{E})_z dx dy = \nabla \wedge \mathbf{E} \cdot d\mathbf{S}.$$

where we have recognised that the term in brackets is the z -component of the vector $\nabla \wedge \mathbf{E}$. The last expression follows because the area vector representing the loop is given by $d\mathbf{S} = \hat{z} dx dy$.

Although we have not proved it, this result is general, *i.e.*, for an infinitesimal flat element of area $d\mathbf{S}$ of any shape and orientation bounded by a loop $C = \partial S$ we can write

$$\oint_{C=\partial S} \mathbf{E} \cdot d\mathbf{l} = \nabla \wedge \mathbf{E} \cdot d\mathbf{S}. \quad (4.2)$$

The quantity $\nabla \wedge \mathbf{E}$ is a *vector* and is called the *curl* of the electric field. You may also sometimes see it called the *rot* \mathbf{E} , short for rotation. Eq. 4.2 defines curl in the same way that Eq. 3.3 defined the divergence.

Of all three derivatives we have now encountered — grad, div, and curl — the curl is the most difficult to get a feel for. Its nature is defined by Eq. 4.2. When thinking about curl, one should picture a small loop embedded in the vector field and consider what the circulation around it is. Still, it is not always obvious whether there is any overall line integral. We will look at some examples later which may help, but first we will finish with Stokes' Theorem.

Any finite surface can be subdivided into many small flat facets obeying the above equation. Adding the line integrals of all of these facets, the individual contributions cancel except on the outer boundary (for example refer back to Fig. 4.1 and consider adding the integrals around the two adjacent dashed squares). We then obtain *Stokes' Theorem*

$$\oint_{C=\partial S} \mathbf{E} \cdot d\mathbf{l} = \int_S \nabla \wedge \mathbf{E} \cdot d\mathbf{S}, \quad (4.3)$$

for *any* surface S bounded by the circuit $C = \partial S$. This applies to *any* physical vector field, not just \mathbf{E} . S and C here are now finite in contrast to Eq. 4.2 and S no longer has to be flat. With Stokes' theorem we can transform line integrals \iff surface integrals.

As a simple application, let us revisit the condition $\oint_C \mathbf{E} \cdot d\mathbf{l} = 0$ which we derived for electrostatic fields in Chapter 2. Since this applies for any circuit, Stokes' theorem implies that

$$\nabla \wedge \mathbf{E} = \mathbf{0},$$

for electrostatic fields ($\mathbf{0}$ is a zero vector). Such a field is said to be *curl-free* or *irrotational*. The study of irrotational fluid flows for which $\nabla \wedge \mathbf{v} = \mathbf{0}$ is of great importance in aerodynamics, and approximations based on this explain why aircraft fly.

The reverse of the above condition is also true. That is if a vector field \mathbf{A} satisfies $\nabla \wedge \mathbf{A} = \mathbf{0}$, then we can write $\oint_C \mathbf{A} \cdot d\mathbf{l} = 0$ and, from Chapter 2, that \mathbf{A} can be derived from a potential $\mathbf{A} = \nabla \phi$. We are saying then that $\nabla \wedge \nabla \phi = \mathbf{0}$ which is reasonable if you remember that the cross-product of a vector with itself is zero (although this is not a proof because ∇ is not an ordinary vector).

Since the condition $\nabla \wedge \mathbf{v} = \mathbf{0}$ implies $\mathbf{v} = \nabla \phi$, curl-free flows are also called *potential flows*. Recalling that incompressible flows satisfy $\nabla \cdot \mathbf{v} = 0$, then we have $\nabla^2 \phi = 0$, and so incompressible potential flows satisfy Laplace's equation which is also satisfied by electrostatic fields, a useful mathematical similarity between very different physical systems.

4.4 Differential version of Faraday's Law

Applying Stokes' theorem to the left-hand side of Faraday's Law we obtain

$$\int_S \nabla \wedge \mathbf{E} \cdot d\mathbf{S} = -\frac{d}{dt} \int_S \mathbf{B} \cdot d\mathbf{S}.$$

Since this applies to *any* loop, we must have

$$\nabla \wedge \mathbf{E} = -\frac{d\mathbf{B}}{dt}.$$

The *total* time derivative d/dt allows the loop to move, but we do not want this because this means that we are measuring \mathbf{B} in our rest frame while we are measuring \mathbf{E} in the rest frame of the loop. A simple thought experiment shows that \mathbf{E} and \mathbf{B} change according to the frame in which they are measured: Picture a charge q moving at velocity \mathbf{v} in a region with a magnetic field \mathbf{B} but no electric field. The force on the charge is $q\mathbf{v} \wedge \mathbf{B}$. How does the picture alter when viewed from a frame in which the charge is at rest (even if it only at rest for an infinitesimal time). Since in this frame the charge is at rest, there is no $\mathbf{v} \wedge \mathbf{B}$ term, and yet the charge must feel a force because it moves in a circle in a magnetic field. We are forced to the conclusion that in the new frame there is an electric field, which for low \mathbf{v} must have strength $\mathbf{v} \wedge \mathbf{B}$. In other words, a magnetic field in one frame may look like an electric field in another frame (see the section on Lorentz transformations below).

The important point is that it is essential to measure electric and magnetic fields in the *same* reference frame. This can be done by fixing the loop to be stationary in our rest frame (and replacing total derivatives by partial derivatives) whereby we obtain

$$\nabla \wedge \mathbf{E} = -\frac{\partial \mathbf{B}}{\partial t} \tag{4.4}$$

This is the differential version of Faraday's Law and the second of Maxwell's equations. It contains no more physics than the integral version, Eq. 4.1, but it applies at a point.

4.4.1 Lorentz Transformations for an Electromagnetic Field

A full description of the transformation of electric and magnetic fields from one frame to another requires the Special Theory of Relativity (and indeed was the main subject of Einstein's original paper) and is beyond the scope of this course (*i.e.*, this subsection is *not* examinable).

For the components of vectors parallel (\parallel) and perpendicular (\perp) to \mathbf{v} , the vector form of the electromagnetic (Lorentz) transformation formulas is

$$\begin{aligned} \mathbf{E}_{\parallel} &= \mathbf{E}'_{\parallel}; & \mathbf{E}_{\perp} &= \gamma(\mathbf{E}'_{\perp} - (\mathbf{v} \wedge \mathbf{B}')_{\perp}); \\ \mathbf{B}_{\parallel} &= \mathbf{B}'_{\parallel}; & \mathbf{B}_{\perp} &= \gamma\left(\mathbf{B}'_{\perp} + \frac{1}{c^2}(\mathbf{v} \wedge \mathbf{E}')_{\perp}\right); \end{aligned}$$

where

$$\gamma = \left(1 - \frac{v^2}{c^2}\right)^{-1/2}.$$

The primed expressions (\mathbf{E}' , \mathbf{B}') correspond to quantities measured in a coordinate system moving at a uniform velocity \mathbf{v} with respect to the coordinate system in which the unprimed expressions, (\mathbf{E} , \mathbf{B}), are deduced. Can you see where the Lorentz force law comes from? What about the Biot-Savart law?

4.5 Calculating the curl

As with the divergence, the direct application of $\text{curl } \mathbf{A} \equiv \nabla \wedge \mathbf{A}$ will in some cases be the simplest approach, although the cross-product can make the calculation of curl error-prone.

- △ The curl of a *vector function* \mathbf{A} is itself a *vector* $\nabla \wedge \mathbf{A}$. You cannot compute the curl of a scalar: that is meaningless. In general, the value of the curl depends on the point at which $\nabla \wedge \mathbf{A}$ is evaluated.

Geometrical interpretation: The name *curl* should indicate to you that $\nabla \wedge \mathbf{A}$ measures how much the vector \mathbf{A} curls around the point in question.

Example 4.2: What is the curl of a field given by $\mathbf{E} = \mathbf{A} \wedge \mathbf{r}$ where \mathbf{A} is a constant vector?

Expanding out the cross-product, the field is given by

$$\begin{pmatrix} E_x \\ E_y \\ E_z \end{pmatrix} = \begin{pmatrix} A_y z - A_z y \\ A_z x - A_x z \\ A_x y - A_y x \end{pmatrix}.$$

We then apply

$$\nabla \wedge \mathbf{E} = \begin{pmatrix} \partial E_z / \partial y - \partial E_y / \partial z \\ \partial E_x / \partial z - \partial E_z / \partial x \\ \partial E_y / \partial x - \partial E_x / \partial y \end{pmatrix} = \begin{pmatrix} 2 A_x \\ 2 A_y \\ 2 A_z \end{pmatrix} = 2 \mathbf{A}.$$

The derivatives are all straightforward, and we obtain the simple result $\nabla \wedge \mathbf{E} = 2 \mathbf{A}$.

Since \mathbf{A} is a constant, $\mathbf{E} = \mathbf{A} \wedge \mathbf{r}$ is the form of field that arises when the magnetic flux has a fixed direction and changes at a constant rate. The electric field in a solenoid has this form.

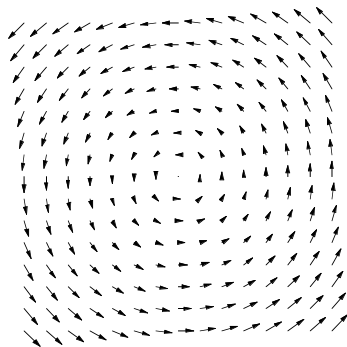
$$\mathbf{A} = \{a, b, c\}; \quad \mathbf{r} = \{x, y, z\};$$

$$\mathbf{E} = \mathbf{A} \wedge \mathbf{r}$$

$$\{bz - cy, cx - az, ay - bx\}$$

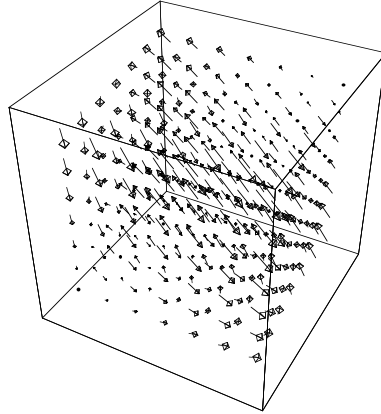
It is usual to visualize this vector field in two dimensions. For example,

<< Graphics`PlotField`; PlotVectorField({-y, x}, {x, -1, 1}, {y, -1, 1});



From the geometrical interpretation we expect this field to have non-zero curl at any point $\{x, y, z\}$. Also, we expect the curl to point in the z -direction as the right-hand rule would suggest. Note that the vector field really exists in three dimensions:

```
PlotVectorField3D({z - y, x - z, y - x}, {x, -1, 1}, {y, -1, 1}, {z, -1, 1}, VectorHeads -> True);
```



However, it is harder to decide from this plot whether we expect this field to have non-zero curl.

Computing the curl in cartesian coordinates we obtain

$$\{\partial_y \mathcal{E}[\mathbf{3}] - \partial_z \mathcal{E}[\mathbf{2}], \partial_z \mathcal{E}[\mathbf{1}] - \partial_x \mathcal{E}[\mathbf{3}], \partial_x \mathcal{E}[\mathbf{2}] - \partial_y \mathcal{E}[\mathbf{1}]\}$$

$$\{2a, 2b, 2c\}$$

```
% == 2 A
True
```

On other occasions it is better to remember the meaning of curl. For example what is the curl of $\mathbf{E} = r^A e^{-r/\lambda} \hat{\mathbf{r}}$? It would be a difficult task to calculate all the necessary derivatives, but also an unnecessary one. This is a spherically symmetric field, so the curl, which is a vector, can only point along the radial direction: anything else would not be spherically symmetric. The derivation of the curl (see Eq. 4.2) means that, in the radial direction we need to calculate the line integral around a loop perpendicular to the radial direction. But this is everywhere perpendicular to the field and therefore the line integral, and the curl, are zero everywhere. This is true for *any* spherically symmetric field, and seems natural as there is no sense of rotation about such a field.

When calculating the divergence we used its basic definition to compute its value in a spherically symmetric case (see section 3.4.1). A similar approach can be taken with the curl in the case of cylindrical symmetry. Consider a field \mathbf{A} that runs in circles around the z -axis with a strength that varies as $A(r)$ where r is the distance from the z -axis. This describes, for instance, the magnetic field around a wire carrying a current, with $A(r) \propto 1/r$.

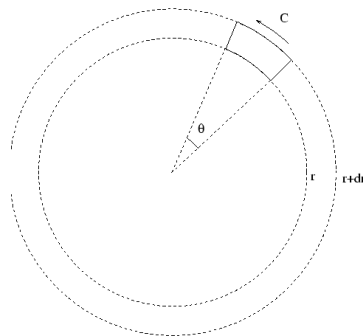


Figure 4.4 The figure shows the path of integration used to evaluate the curl in a field that runs in circles around an axis pointing out of the page and passing through the centre of the circles.

Only loops perpendicular to z can have any line integral. Therefore, the curl must lie in the z -direction. Therefore we choose a small loop which lies perpendicular to the z -axis as illustrated in Fig. 4.4. We are free to choose any shape for the loop, so we pick one to make the calculation as easy as possible. The circuit C illustrated in Fig. 4.4 either runs parallel to the field or perpendicular to it. The line integral around it can then be written down as

$$\oint_C \mathbf{A} \cdot d\mathbf{l} \equiv \nabla \wedge \mathbf{A} \cdot d\mathbf{S} = (r + dr) A(r + dr) \theta - r A(r) \theta = \frac{\partial(r A(r))}{\partial r} \theta dr = |\nabla \wedge \mathbf{A}| r \theta dr$$

with the last step following from Eq. 4.2, and where $A(r)$ is the magnitude of the field at distance r . We can easily verify the middle result (up to second order) using *Mathematica*:

$$(r + dr) A(r + dr) \theta - r A(r) \theta + O(dr)^2 = \frac{\partial(r A(r))}{\partial r} \theta dr + O(dr)^2$$

True

Therefore we obtain

$$\nabla \wedge \mathbf{A} = \frac{1}{r} \frac{\partial(r A(r))}{\partial r} \hat{z}.$$

Proceeding in this manner one can obtain more general expressions for the curls of fields expressed in any coordinate system. See Appendix A for a summary of vector operators in orthogonal coordinate systems. For example, the general expression for the curl in spherical polar coordinates is (see Eq. A.16)

$$\begin{aligned} \nabla \wedge \mathbf{A} = & \frac{1}{r \sin(\theta)} \left(\frac{\partial(\sin(\theta) A_\phi)}{\partial \theta} - \frac{\partial A_\theta}{\partial \phi} \right) \mathbf{e}_r + \\ & \frac{1}{r \sin(\theta)} \left(\frac{\partial A_r}{\partial \phi} - \sin(\theta) \frac{\partial(r A_\phi)}{\partial r} \right) \mathbf{e}_\theta + \frac{1}{r} \left(\frac{\partial(r A_\theta)}{\partial r} - \frac{\partial A_r}{\partial \theta} \right) \mathbf{e}_\phi \end{aligned}$$

where $\mathbf{A} \equiv A_r \hat{r} + A_\theta \hat{\theta} + A_\phi \hat{\phi}$.

Example 4.3: What is the curl of a field that runs in circles around the z -axis with a strength that drops off inversely with distance?

Using the above result, we put $A(r) = 1/r$ and find that $\nabla \wedge \mathbf{A} = \mathbf{0}$.

This is the form of magnetic field near a wire carrying a current. It illustrates the point that a field can appear to have circulation but have no curl. However our calculation breaks down on the axis itself because a loop enclosing that would definitely have a finite integral (*c.f.* computing the divergence of the electric field of a point charge).

Example 4.4: What is the curl of a field that runs in circles around the z -axis with a strength that increases linearly with distance?

Now $A(r) = r$ and we find $\nabla \wedge \mathbf{A} = 2 \hat{z}$.

This is the form of magnetic field inside a wire of uniform current density. It turns out that the curl of the magnetic field is proportional to the current density so it is no fluke that the curl turns out to have constant magnitude. Since the region near a current carrying wire carries no current itself, the result in Example 4.3 is no surprise either.

4.6 General remarks on grad, div and curl

The curl is the last new vector derivative. It is important to appreciate the nature of these quantities in the sense that $\text{grad } \phi = \nabla \phi$ is a vector, as is $\text{curl } \mathbf{E} = \nabla \wedge \mathbf{E}$, whereas $\text{div } \mathbf{E} = \nabla \cdot \mathbf{E}$ is a scalar. Realising this helps one avoid mistakes of the following kind

$$\nabla \cdot \mathbf{E} = -\frac{\partial B}{\partial t}$$

because a scalar on the left cannot equal a vector on the right. When first met these quantities can be confusing. One has no "feel" or "intuition" for them. Intuition is actually a misleading expression; "experience of" would be more accurate. Why should anyone have intuition for a concept such as curl which they have never met before? You can only develop "intuition" after use and after seeing these quantities in action. The way to develop it most quickly is to remember equations such as Eq. 4.4 and always to focus on the physical meaning behind the symbols.

4.7 Summary

We started in this chapter by expressing Faraday's law. We then considered how this can be applied to an infinitesimal region and in doing so defined a new quantity, the curl of \mathbf{E} , defined by (for infinitesimal loops)

$$\oint_C \mathbf{E} \cdot d\mathbf{l} \equiv \text{curl } \mathbf{E} \cdot d\mathbf{S} = \nabla \wedge \mathbf{E} \cdot d\mathbf{S},$$

with the second form based on consideration of a small rectangular loop. The curl of a vector field is itself a vector.

This led on to more general equation called *Stokes' Theorem* that can be applied to finite surfaces and can be used to transform between surface and line integrals. This equation was then applied to the integral version of Faraday's law to arrive at the differential form of Faraday's law.

Chapter 5

Magnetic fields

5.1 Introduction

In this chapter we look at the physics of magnetostatics. We will encounter the second pair of Maxwell's equations although one of them will have to be modified for time variable phenomena later. We make use of Gauss' and Stokes' theorems, but no new mathematics has to be introduced. Our starting points are the Biot-Savart and Ampère's laws.

5.2 The Biot-Savart Law

The Biot-Savart law gives the contribution to the magnetic at a point from a small current element. Let a current I flow through a short element $d\mathbf{l}$. Then the magnetic field due to this element at a point r away from it is given by

$$d\mathbf{B} = \frac{\mu_0}{4\pi} \frac{I d\mathbf{l} \wedge \hat{\mathbf{r}}}{r^2}$$

The cross-product here gives a magnetic field obeying the usual right-hand rule for magnetic fields from currents. That is, with the thumb of your right-hand pointing along the current, your fingers point in the direction of the field.

There are rather few situations where the Biot-Savart law proves practical to use and we are not going to use it to any great extent. The main point we take from it is that the field lines run in circles around an axis defined by the direction of $d\mathbf{l}$. If they run in circles, no flux of \mathbf{B} is produced or destroyed, and therefore we can write immediately

$$\nabla \cdot \mathbf{B} = 0, \tag{5.1}$$

because $\nabla \cdot \mathbf{B}$, the divergence of \mathbf{B} , represents the amount of magnetic flux produced per unit volume. This equation is the third of Maxwell's equations. It can be proved more formally, but the proof is not illuminating. If we compare with the equivalent equation for the electric field (Eq. 3.5) $\nabla \cdot \mathbf{E} = \frac{\rho}{\epsilon_0}$, we can interpret $\nabla \cdot \mathbf{B} = 0$ as saying that there are no sources of magnetic flux, or in other words there are no magnetic charges (*magnetic monopoles*).

5.2.1 Lorentz Transformation

As an aside, a current element $I d\mathbf{l}$ can be thought of as a charge moving with a velocity, *i.e.*, $I d\mathbf{l} \equiv q \mathbf{v}$. Hence the Biot-Savart law says that the magnetic field due to this charge at a point r away from it is given by

$$\mathbf{B} = \frac{\mu_0}{4\pi} \frac{q \mathbf{v} \wedge \hat{\mathbf{r}}}{r^2}.$$

We can relate this to the electric field of the point charge as follows

$$\mathbf{B} = \frac{\mu_0}{4\pi} 4\pi \epsilon_0 \mathbf{v} \wedge \left(\frac{1}{4\pi \epsilon_0} \frac{\hat{\mathbf{r}}}{r^2} \right) = \mu_0 \epsilon_0 \mathbf{v} \wedge \mathbf{E}.$$

Compare this with the Lorentz transformation formula for magnetic fields:

$$\mathbf{B}_\perp = \gamma \left(\mathbf{B}'_\perp + \frac{1}{c^2} (\mathbf{v} \wedge \mathbf{E}')_\perp \right);$$

5.3 Ampère's Circuital Law

Ampère's circuital law (often just Ampère's law) is not independent from the Biot-Savart law, and can be derived from it. In fact as Ampère did most of the experiments that lead up to both laws, there is a case for renaming the Biot-Savart law, except for the confusion it would cause. The actual derivation is again not very informative, and we are content to quote the result, which in words says that the line integral of \mathbf{B} around a closed loop is equal to μ_0 times the total current through the loop.

Following the work of previous chapters, we can immediately write this in symbolic form as

$$\oint_{C=\partial S} \mathbf{B} \cdot d\mathbf{l} = \mu_0 \int_S \mathbf{J} \cdot d\mathbf{S}. \quad (5.2)$$

Here C is the loop through which the current flows, \mathbf{J} is the current density and S is any surface bounded by C . Applying Stokes' theorem to transform the line integral into a surface integral we have:

$$\oint_{C=\partial S} \mathbf{B} \cdot d\mathbf{l} = \int_S \nabla \wedge \mathbf{B} \cdot d\mathbf{S} = \mu_0 \int_S \mathbf{J} \cdot d\mathbf{S}$$

and, since this applies for any loop C , we must have

$$\nabla \wedge \mathbf{B} = \mu_0 \mathbf{J}, \quad (5.3)$$

which is the differential form of Ampère's law. This was the relation referred to in the discussion after example 4.4 in which we calculated the curl of some example fields.

As an unusual application of Ampère's law, suppose that we wish to measure the total current flowing to or from the ground during a thunderstorm. We could do so by measuring the magnetic field at a series of points on the ground at the boundary of the storm. Taking the line integral would give us the current. This would be a great deal easier than measuring the current directly, which would in any case require knowing where lightning was going to strike.

Example 5.1: A current I flows in a long wire of circular cross-section of radius a (Fig. 5.1). What is the magnetic field as a function of the distance r from the axis of the wire?

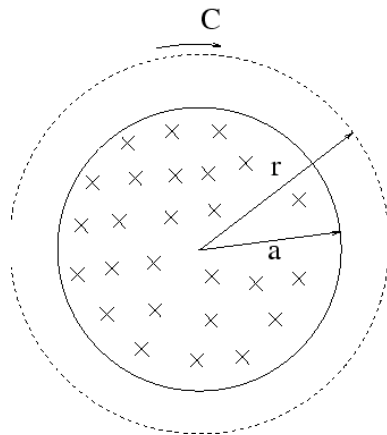


Figure 5.1 A cross-section of a wire carrying a current into the page (represented by crosses). C is the path used to determine the magnetic field.

We will apply the integral form of Ampère's law (Eq. 5.2).

We need to define a suitable circuit. Since the magnetic field must run around the wire in circles, the obvious path is itself a circle centred on the axis of the wire so that the magnetic field is everywhere parallel to it and of the same strength. This problem is very similar to example 4.1 where we calculated the electric field of a solenoid. Fig. 5.1 shows such a path.

The line integral $\oint_C \mathbf{B} \cdot d\mathbf{l}$ reduces to $2\pi r B$. The current linked depends on whether the circuit is inside or outside the wire. If it is outside ($r > a$) then the current enclosed is simply I ; if it is inside ($r < a$) then the current enclosed scales with area (*i.e.*, r^2) and must therefore be $I(r/a)^2$ so that it equals I for $r = a$. Applying Ampère's law (Eq. 5.2) we obtain

$$B = \begin{cases} \frac{\mu_0 I}{2\pi r}, & r > a \\ \frac{\mu_0 I r}{2\pi a^2}, & r < a \end{cases}.$$

Example 5.2: Derive the equation $B = \mu_0 n I$ for the magnetic field inside a long solenoid with n turns per unit length and carrying a current I .

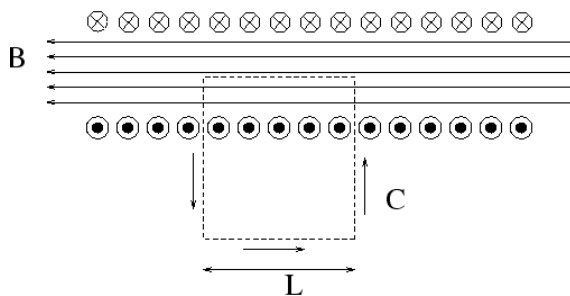


Figure 5.2 Figure 5.2: Cross-section of a solenoid with \otimes representing wires carrying current into the page and \odot representing currents flowing out of the page. The circuit C is used to determine the magnetic field.

Consider the rectangular circuit as shown in Fig. 5.2. Only the side running parallel to the field inside the solenoid gives any contribution to $\oint_C \mathbf{B} \cdot d\mathbf{l}$. Its contribution is BL . The circuit links a current of nLI and so $BL = \mu_0 nLI$. Therefore $B = \mu_0 nI$.

5.4 Summary

Starting from the Biot-Savart law, the third of Maxwell's equations, Eq. 5.1, was written down. This equation expresses the fact that no magnetic charges have ever been found. Next Ampère's law was translated into mathematical form (Eq. 5.2). Applying Stokes' theorem, the differential version, Eq. 5.3, was immediately obtained.

Chapter 6

Electromagnetic Waves

6.1 Introduction

In this chapter we show that Ampère's Law cannot apply in the time-varying case. We consider how to modify, introducing an extra term, the *displacement current*, to satisfy charge conservation. We then demonstrate the existence of electromagnetic waves. We examine the essential properties of these waves in the vacuum, considering both the general properties of waves and properties specific to electromagnetic waves.

6.2 The displacement current

In the previous chapters we have derived the following differential equations describing electric and magnetic fields

$$\begin{aligned}\nabla \cdot \mathbf{E} &= \frac{\rho}{\epsilon_0} & \nabla \cdot \mathbf{B} &= 0 \\ \nabla \wedge \mathbf{E} &= -\frac{\partial \mathbf{B}}{\partial t} & \nabla \wedge \mathbf{B} &= \mu_0 \mathbf{J}\end{aligned}$$

The second pair of equations relate the curl of one vector field to a different vector field. If we take the final equation, for instance, it says that the free current density, \mathbf{J} , is the curl of the magnetic field. This places an important restriction upon the nature of \mathbf{J} . To realise why, we first need a mathematical result (a vector field identity) which states that for any vector field \mathbf{A} , the divergence of its curl equals zero, *i.e.*,

$$\nabla \cdot (\nabla \wedge \mathbf{A}) = 0.$$

Taking the divergence ($\nabla \cdot$) of both sides of $\nabla \wedge \mathbf{B} = \mu_0 \mathbf{J}$ we obtain $\nabla \cdot \mathbf{J} = 0$. In other words Ampère's Law (as we have seen it so far) implies that the current density \mathbf{J} is divergence-less.

This result cannot be true. It says that the total flux of current per unit volume is everywhere zero. Equivalently, using Gauss' Theorem we have

$$\int_V \nabla \cdot \mathbf{J} \, d\mathbf{r} = \oint_{S=\partial V} \mathbf{J} \cdot d\mathbf{S} = 0,$$

which says that the total current flowing out of any volume is always zero. This is wrong because it would mean that nothing could ever be charged or discharged. Every time a capacitor is charged, $\nabla \cdot \mathbf{J} = 0$ is violated as charge flows on and off the plates.

We saw the correct relation for the divergence of the current density when we discussed continuity equations. If current flows out of a volume, it is balanced by a loss of charge from the volume, and this led us to $\nabla \cdot \mathbf{J} + \frac{\partial \rho}{\partial t} = 0$.

Using Gauss' Law $\nabla \cdot \mathbf{E} = \frac{\rho}{\epsilon_0}$ we may write the continuity equation as

$$\begin{aligned}\frac{\partial \rho}{\partial t} &= \epsilon_0 \frac{\partial \nabla \cdot \mathbf{E}}{\partial t} = \epsilon_0 \nabla \cdot \frac{\partial \mathbf{E}}{\partial t} \\ \nabla \cdot \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right) &= 0,\end{aligned}$$

where we have made use of the commutativity of ∇ and $\frac{\partial}{\partial t}$, which is to say that the order in which they are applied makes no difference. Therefore if Ampère's Law is changed to

$$\nabla \wedge \mathbf{B} = \mu_0 \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right), \quad (6.1)$$

we have a divergence-less field on each side of the equation. Moreover, in static cases the modified equation reduces to $\nabla \wedge \mathbf{B} = \mu_0 \mathbf{J}$ which was derived from magnetostatic experiments. Finally, the modified equation now resembles Faraday's Law, the only difference being that because there are no magnetic charges, there is no magnetic current term in Faraday's Law. Equation 6.1 is our final version of Ampère's Law and completes the set of Maxwell's equations.

The new term, $\mu_0 \epsilon_0 \frac{\partial \mathbf{E}}{\partial t}$, is related to the *displacement current* (actually a current density) and was introduced by Maxwell. Although very suggestive, the above discussion provides only a motivation for the introduction of the displacement current and ultimately its true test rests on experiments. For example, there are many other terms that could be added which would also have zero divergence. However, the displacement current is needed for the propagation of electromagnetic waves, and so every time one turns on a light its existence is demonstrated.

The effects of the displacement current are exactly those of an ordinary current, and cannot be distinguished from it. For instance, as a capacitor is charged, and the field between the plates increases, it is as if a current were flowing between the plates and a magnetic field will be generated just as it would if there were a true current of the same magnitude.

Why wasn't the displacement current found experimentally? First, the experiments that led to Ampère's Law are difficult to perform in time varying cases; Ampère experimented with coils and steady currents. Second the displacement current is small. In vacuum, a rate of change of electric field of order $10^{11} \text{ V.m}^{-1}.\text{s}^{-1}$ over 1 m^2 is needed to produce a current of only 1 A. The displacement current is often negligible with the important exception of when no ordinary current can flow, as in a vacuum or a dielectric. When we derive the wave equation in Section 6.5, the displacement current is vital.

6.3 Summary of Maxwell's Equations

Maxwell's equations, in terms of total charge and current densities, ρ and \mathbf{J} , read

$$\begin{aligned}\nabla \cdot \mathbf{E} &= \frac{\rho}{\epsilon_0} & \nabla \cdot \mathbf{B} &= 0 \\ \nabla \wedge \mathbf{E} &= -\frac{\partial \mathbf{B}}{\partial t} & \nabla \wedge \mathbf{B} &= \mu_0 \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right)\end{aligned} \quad (6.2)$$

Note that these equations apply generally but, because the total charge and current densities include contributions from polarisation and magnetisation, it is not usually convenient to use them when materials are present.

Each of these differential equations has an integral equivalent. Very often it is the integral versions that are easier to apply, but the differential equations are vital in the study of electromagnetic radiation as we will see in the rest of this chapter. The integral versions can be derived by suitable integration followed by application of Stokes' Theorem or Gauss' Theorem. For example, consider

$$\nabla \wedge \mathbf{B} = \mu_0 \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right).$$

This has a curl on the left, so if we integrate it over some surface, we will be able to transform the resulting surface integral into a line integral by Stokes' theorem. Thus we get the following steps

$$\int_S \nabla \wedge \mathbf{B} \cdot d\mathbf{S} = \mu_0 \int_S \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right) \cdot d\mathbf{S},$$

and so, after applying Stokes' Theorem to the left-hand side, we get

$$\oint_{C=\partial S} \mathbf{B} \cdot d\mathbf{l} = \mu_0 \int_S \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right) \cdot d\mathbf{S},$$

This equation says that the line integral of \mathbf{B} around a circuit is equal to the sum of the free and displacement currents flowing through it.

We can go through a similar procedure for each equation and we obtain the following integral equations equivalent to Equations 6.2:

$$\begin{aligned} \oint_S \mathbf{E} \cdot d\mathbf{S} &= \frac{1}{\epsilon_0} \int_V \rho \, d\mathbf{r} & \oint_S \mathbf{B} \cdot d\mathbf{S} &= 0 \\ \oint_C \mathbf{E} \cdot d\mathbf{l} &= - \int_S \frac{\partial \mathbf{B}}{\partial t} \cdot d\mathbf{S} & \oint_{C=\partial S} \mathbf{B} \cdot d\mathbf{l} &= \mu_0 \int_S \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right) \cdot d\mathbf{S} \end{aligned} \quad (6.3)$$

The right-hand sides of these equations can be replaced by integrated quantities such as charge or current as appropriate.

6.4 General properties of waves

6.4.1 Phase or wave velocity

The 3D wave equation has the form

$$\nabla^2 \zeta = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta}{\partial t^2},$$

where ζ represents any wave-like quantity. Why does this describe waves? First we have to define what a wave is. A wave is some sort of disturbance that propagates with time. In the simplest case waves propagate without changing shape. For example, someone's voice sounds the same, apart from loudness, almost independently of the distance of the speaker. The sound waves are little distorted by travel in air. A wave of this form travelling in the x -direction can be described by

$$\zeta(x, t) = \zeta_0 f(x - vt).$$

This function is constant for constant values of $x - vt$, which implies that $x = vt + \text{const}$ and so v represents the speed at which the disturbance travels which we will call the *wave* or *phase velocity*.

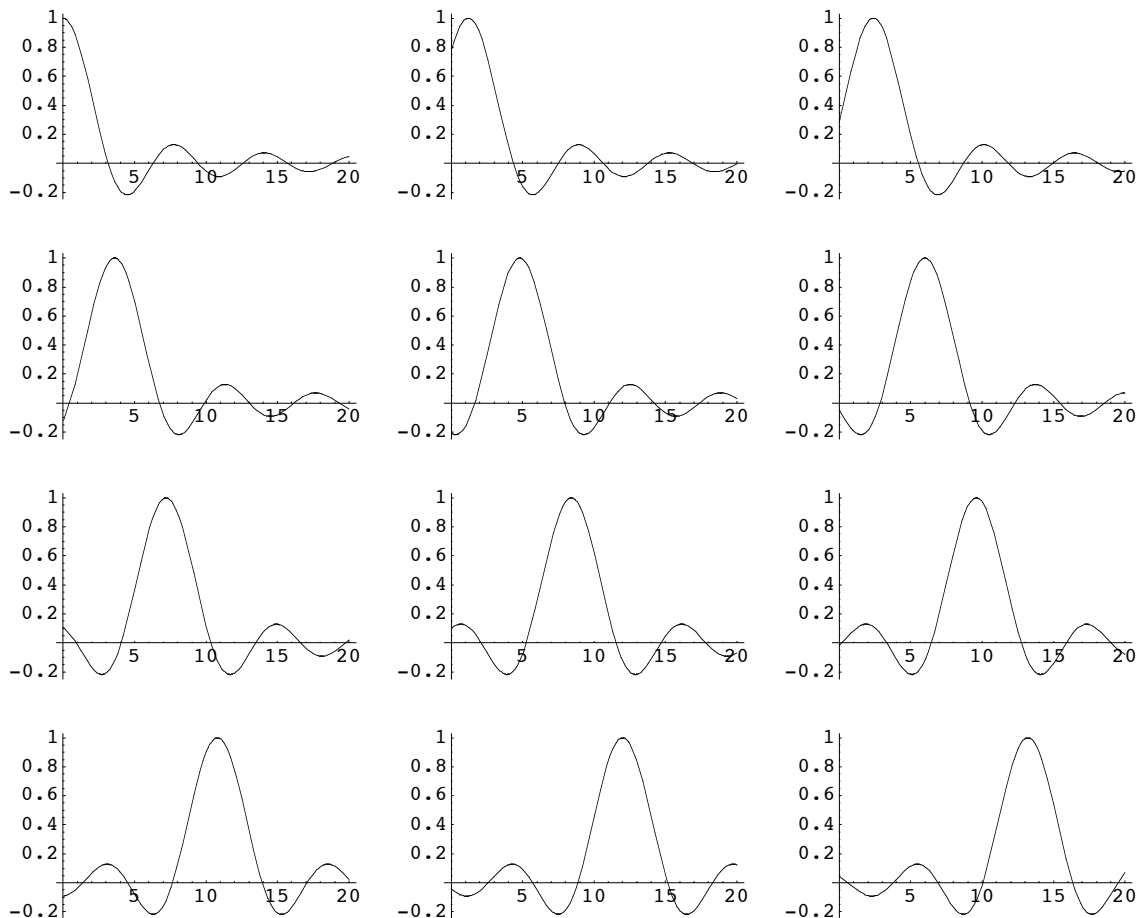
For example, with the function

$$f(x) = \frac{\sin(x)}{x};$$

double-click on the graphic below to see an animation of a one-dimensional wave.

Table[Plot[f[x - 1.2 t], {x, 0.01, 20}, PlotRange -> All], {t, 0, 11}];

Show[GraphicsArray[Partition[%, 3]]];



To see that $\zeta(x, t)$ satisfies the wave equation, we substitute it in. We therefore need to calculate various derivatives. Thus, for example, putting

$$\zeta(x, t) = \zeta_0 f(x - vt);$$

then the second derivative with respect to x is

$$\frac{\partial^2 \zeta(x, t)}{\partial x^2} = \zeta_0 f''(x - vt)$$

whilst the second derivative with respect to t is

$$\frac{\partial^2 \zeta(x, t)}{\partial t^2} = v^2 \zeta_0 f''(x - vt)$$

Substituting into the wave equation we find

$$\frac{\partial^2 \zeta(x, t)}{\partial x^2} = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta(x, t)}{\partial t^2}$$

$$\zeta_0 f''(x - tv) = \frac{v^2 \zeta_0 f''(x - tv)}{v_\phi^2}$$

Since $f''(x - tv)$ crops up on both sides, this result holds for *any* f , provided that

$$v^2 = v_\phi^2,$$

so that the constant v_ϕ appearing in wave equation can be identified with the wave velocity. The above relation shows that $v = \pm v_\phi$ because waves can go in either direction.

6.4.2 Linearity and superposition

The next important property of the wave equation is its *linearity* which allows one to superpose solutions. This means that given any two solutions of the wave equation, their sum (or in general any linear combination) is also a solution. The physical consequence of this property is that two beams of light do not affect each other even where they cross. Linearity is an extremely useful property, and although nonlinear equations show more interesting effects, they are often harder to deal with.

Sound waves are linear at typical strengths; when they become non-linear they turn into shock waves. Ocean waves are approximately linear when the depth of the water is much larger than their height. However, as they approach the shore this is no longer the case and the top of the wave curls over and the wave breaks. The behaviour in this zone is highly nonlinear.

To prove the property of superposition for the wave equation, suppose that we have two solutions ζ_1 and ζ_2 that satisfy the wave equation. That is

$$\nabla^2 \zeta_1 = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta_1}{\partial t^2}, \quad \nabla^2 \zeta_2 = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta_2}{\partial t^2}.$$

Then $\zeta_3 = \zeta_1 + \zeta_2$ is also a solution:

$$\nabla^2 \zeta_3 = \nabla^2 (\zeta_1 + \zeta_2) = \nabla^2 \zeta_1 + \nabla^2 \zeta_2 = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta_1}{\partial t^2} + \frac{1}{v_\phi^2} \frac{\partial^2 \zeta_2}{\partial t^2} = \frac{1}{v_\phi^2} \frac{\partial^2 (\zeta_1 + \zeta_2)}{\partial t^2} = \frac{1}{v_\phi^2} \frac{\partial^2 \zeta_3}{\partial t^2}.$$

If there were any non-linear terms such as ζ^2 , the above proof would break down. Hence the close connection between linearity and superposition.

6.4.3 Plane waves

We saw above that $\zeta(x, t) = \zeta_0 f(x - v_\phi t)$ is a solution of the wave equation corresponding to a wave travelling at v_ϕ in the x -direction. More generally

$$\zeta(\mathbf{r}, t) = \zeta_0 f(\mathbf{k} \cdot \mathbf{r} - \omega t)$$

is a solution. To confirm that this is a possible solution, we write this equation out explicitly

$$\zeta(\mathbf{x}_-, \mathbf{y}_-, \mathbf{z}_-, \mathbf{t}_-) = \zeta_0 f(a x + b y + c z - \omega t);$$

and compute the left-hand side

$$\frac{\partial^2 \zeta(x, y, z, t)}{\partial x^2} + \frac{\partial^2 \zeta(x, y, z, t)}{\partial y^2} + \frac{\partial^2 \zeta(x, y, z, t)}{\partial z^2} \quad // \text{Factor}$$

$$(a^2 + b^2 + c^2) \zeta_0 f''(a x + b y + c z - t \omega)$$

and right-hand side

$$\frac{1}{v_\phi^2} \frac{\partial^2 \zeta(x, y, z, t)}{\partial t^2}$$

$$\frac{\omega^2 \zeta_0 f''(a x + b y + c z - t \omega)}{v_\phi^2}$$

of the wave equation. The wave equation is satisfied provided that

$$\frac{\omega^2}{v_\phi^2} = a^2 + b^2 + c^2$$

where a , b , and c are the components of the vector $\mathbf{k} = (a, b, c)$ which has magnitude k . Then we have the important (dispersion) relation

$$v_\phi = \frac{\omega}{k} \quad (6.4)$$

The function $f(a x + b y + c z - \omega t)$ is constant for x , y , z , and t which satisfy

$$a x + b y + c z - \omega t = \text{constant},$$

At a fixed instant of time therefore

$$a x + b y + c z = \mathbf{k} \cdot \mathbf{r} = \text{constant}.$$

This is the equation of a plane in three dimensions and so the function f describes *plane waves*. All physical quantities related to the wave are constant over these planes. The vector \mathbf{k} is perpendicular to the planes of constant phase and therefore is parallel to the direction of the waves. It is called the *wave vector*.

As time changes, the wave-fronts move. We can calculate the velocity at which they move from consideration of the argument of f , more usually called the *phase*, which we denote by $\psi = \mathbf{k} \cdot \mathbf{r} - \omega t$. Differentiating with respect to time, with ψ constant and setting $\mathbf{v} = \frac{d\mathbf{r}}{dt}$ we have $\mathbf{k} \cdot \mathbf{v} = \omega$. Movement parallel to planes of constant phase cannot be seen so it is only movement perpendicular to the wavefronts that is of interest. The wave vector \mathbf{k} is perpendicular to the wave fronts and so we can set $\mathbf{v} = v_\phi \hat{\mathbf{k}}$, whereby we obtain

$$v_\phi = \frac{\omega}{k},$$

which we have already seen when we proved that the plane wave solution satisfied the wave equation. This shows that the v_ϕ in the wave equation is the velocity at which the wavefronts move in the direction of the wave vector \mathbf{k} , as expected. Thus our 1D results extend naturally to 3D through the wave vector.

The wave vector \mathbf{k} points in the direction of motion of the wavefronts and for all the waves that we will consider, this coincides with the direction that a beam of radiation (*i.e.*, the energy) travels. Rather surprisingly perhaps, there are waves for which this is not the case (*e.g.*, light in birefringent crystals) but we will not consider these in this course.

6.4.4 Harmonic plane waves

A particularly important solution to the wave equation is the *harmonic plane wave* form

$$\zeta(\mathbf{r}, t) = \zeta_0 e^{i(\mathbf{k}\cdot\mathbf{r} - \omega t)} = \zeta_0 (\cos(\mathbf{k}\cdot\mathbf{r} - \omega t) + i \sin(\mathbf{k}\cdot\mathbf{r} - \omega t))$$

Applying the operators $\frac{\partial}{\partial t}$ and ∇ to $\zeta(\mathbf{r}, t)$ we see that

$$\begin{aligned} \frac{\partial}{\partial t} \zeta(\mathbf{r}, t) &= -i \omega \zeta(\mathbf{r}, t), \\ \nabla \zeta(\mathbf{r}, t) &= i \mathbf{k} \zeta(\mathbf{r}, t), \end{aligned}$$

along with analogous results for vector fields.

In other words, application of the differential operator $\frac{\partial}{\partial t}$ is equivalent to multiplication by $-i \omega$. Similarly ∇ is equivalent to multiplication by $i \mathbf{k}$. For example, $\nabla \cdot \mathbf{E} = 0$ becomes $i \mathbf{k} \cdot \mathbf{E} = 0$ or $\mathbf{k} \cdot \mathbf{E} = 0$.

△ Compare with the quantum mechanical momentum operator, $\mathbf{p} = -i \nabla$.

6.5 Wave equation from Maxwell's Equations

The displacement current term gives us electromagnetic waves. The second pair of Maxwell's equations connects spatial derivatives of each field to the rate of change of the other. It is this coupling from the electric to the magnetic field and back again that allows the propagation of waves. Qualitatively, a time-varying \mathbf{B} -field in $\nabla \wedge \mathbf{E} = -\frac{\partial \mathbf{B}}{\partial t}$, leads to a time-varying \mathbf{E} -field, which leads (self-consistently) to a time-varying \mathbf{B} -field in $\nabla \wedge \mathbf{B} = \mu_0 (\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t})$.

In this course we will only consider electromagnetic waves in vacuum and restrict attention to the situation where there are no free currents ($\mathbf{J} = \mathbf{0}$) or charges ($\rho = 0$). Hence Maxwell's equations become

$$\begin{aligned} \nabla \cdot \mathbf{E} &= 0 & \nabla \cdot \mathbf{B} &= 0 \\ \nabla \wedge \mathbf{E} &= -\frac{\partial \mathbf{B}}{\partial t} & \nabla \wedge \mathbf{B} &= \mu_0 \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \end{aligned} \quad (6.5)$$

The first pair of equations restricts the variety of possible fields.

The standard way to proceed with such pairs of coupled differential equations is to take the derivative of one of them and use the other to eliminate one or other of the independent variables. In this case we take the *curl* of Faraday's Law because this leads to a $\nabla \wedge \mathbf{E}$ term which can be eliminated using Ampère's law:

$$\nabla \wedge \nabla \wedge \mathbf{B} = \mu_0 \epsilon_0 \nabla \wedge \frac{\partial \mathbf{E}}{\partial t}$$

The left-hand side can be rearranged using a second vector field identity

$$\nabla \wedge \nabla \wedge \mathbf{A} = \nabla(\nabla \cdot \mathbf{A}) - \nabla^2 \mathbf{A}.$$

Using the commutativity of ∇ and $\partial / \partial t$ on the right-hand side we obtain

$$-\nabla^2 \mathbf{B} = \mu_0 \epsilon_0 \frac{\partial(\nabla \wedge \mathbf{E})}{\partial t} = -\mu_0 \epsilon_0 \frac{\partial^2 \mathbf{B}}{\partial t^2}$$

Substituting from Equation (6.6) for $\nabla \cdot \mathbf{E}$ and $\nabla \wedge \mathbf{B}$ we then have

$$\nabla^2 \mathbf{E} = \mu_0 \epsilon_0 \frac{\partial^2 \mathbf{E}}{\partial t^2}.$$

This equation has the form of a three-dimensional (*vector*) wave equation for the field \mathbf{E} , with phase velocity

$$v = \frac{1}{\sqrt{\mu_0 \epsilon_0}}$$

Since $\mu_0 = 4\pi \times 10^{-7}$ and $\epsilon_0 = 8.89 \times 10^{-12}$ we obtain

$$v_\phi = 2.99 \times 10^8 \text{ m.s}^{-1},$$

which is equal to the speed of light, c . The clear implication is that light is itself an electromagnetic wave. This is an early instance of the unification of seemingly separate branches of physics, in this case electromagnetism and optics, and one of the triumphs of 19th century physics.

6.6 Relations between fields and the wave vector

In deriving the wave equation we lose information on the relationships between the \mathbf{E} and \mathbf{B} fields. We need to go back to Equation (6.6) to derive these. Using the harmonic plane wave form for the \mathbf{E} and \mathbf{B} fields

$$\mathbf{E}(\mathbf{r}, t) = \mathbf{E}_0 e^{i(\mathbf{k}\cdot\mathbf{r} - \omega t)}$$

$$\mathbf{B}(\mathbf{r}, t) = \mathbf{B}_0 e^{i(\mathbf{k}\cdot\mathbf{r} - \omega t)}$$

we can simplify the equations, translating them from vector differential equations to plain vector equations:

$$\begin{aligned} \mathbf{k} \cdot \mathbf{E} &= 0 & \mathbf{k} \cdot \mathbf{B} &= 0 \\ \mathbf{k} \wedge \mathbf{E} &= \omega \mathbf{B} & \mathbf{k} \wedge \mathbf{B} &= -\mu_0 \epsilon_0 \omega \mathbf{E} \end{aligned} \quad (6.6)$$

The first two equations show that \mathbf{E} and \mathbf{B} are perpendicular to the wave vector \mathbf{k} . Since \mathbf{k} points in the direction of the wave, this means that electromagnetic waves are *transverse waves*. In general a vector in 3D has three degrees of freedom. The condition that \mathbf{E} must be perpendicular to \mathbf{k} reduces this to two degrees of freedom. Physically this corresponds to the two polarisations that light can be split into.

The other two equations relate \mathbf{E} and \mathbf{B} . It is normal to regard the electric field as the one which defines the wave, and for example the direction it points defines the polarisation of the wave. Thus it is convenient to use $\mathbf{k} \wedge \mathbf{E} = \omega \mathbf{B}$ to obtain the magnetic field strength. This equation shows that \mathbf{B} is perpendicular to \mathbf{E} , and so we have found the property of electromagnetic waves that \mathbf{E} , \mathbf{B} , and \mathbf{k} are mutually perpendicular.

Since \mathbf{k} and \mathbf{E} are perpendicular, in terms of magnitudes we have $kE = \omega B$, and therefore $E = \frac{\omega}{k} B = v_\phi B$ by Eq. 6.5. For waves in a vacuum $v_\phi = c$, and so $B = E/c$. The final equation $kB = \mu_0 \epsilon_0 \omega E$ tells us nothing new since with $\mu_0 \epsilon_0 = \frac{1}{c^2}$ it also reduces to $B = E/c$.

6.7 Summary

In this chapter we showed that Ampère's law $\nabla \wedge \mathbf{B} = \mu_0 \mathbf{J}$ fails to satisfy charge conservation and introduced a new term, the displacement current, in order to correct it. Thus we obtained

$$\nabla \wedge \mathbf{B} = \mu_0 \left(\mathbf{J} + \epsilon_0 \frac{\partial \mathbf{E}}{\partial t} \right).$$

This gave us our final versions of Maxwell's equations.

We then studied general properties of the wave equation, and in particular plane waves of the form

$$\zeta(\mathbf{r}, t) = \zeta_0 f(\mathbf{k} \cdot \mathbf{r} - v t).$$

We showed that \mathbf{k} is a vector pointing in the direction of the wave and that the phase or wave velocity v_ϕ was given by

$$v_\phi = \frac{\omega}{k}$$

We then showed that Maxwell's equations in free space lead to a 3D wave equation

$$\nabla^2 \mathbf{E} = \mu_0 \epsilon_0 \frac{\partial^2 \mathbf{E}}{\partial t^2}.$$

The complex exponential form of the wave allowed us to substitute $-i\mathbf{k}$ for ∇ and $i\omega$ for $\partial/\partial t$ in all Maxwell's equations and therefore derive relations between the fields and \mathbf{k} and we showed that \mathbf{E} , \mathbf{B} , and \mathbf{k} were mutually perpendicular.

6.8 Vector field identities

$$\nabla \cdot \nabla \wedge \mathbf{A} = 0$$

It is easy to prove that

$$\nabla \cdot \nabla \wedge \mathbf{A} = 0,$$

by writing out the terms explicitly:

$$\partial_x (\partial_y A_z(x, y, z) - \partial_z A_y(x, y, z)) + \partial_y (\partial_z A_x(x, y, z) - \partial_x A_z(x, y, z)) + \partial_z (\partial_x A_y(x, y, z) - \partial_y A_x(x, y, z)) = 0$$

$$\nabla \wedge \nabla \wedge \mathbf{A} = \nabla(\nabla \cdot \mathbf{A}) - \nabla^2 \mathbf{A}$$

There are elegant methods of proving that

$$\nabla \wedge \nabla \wedge \mathbf{A} = \nabla(\nabla \cdot \mathbf{A}) - \nabla^2 \mathbf{A}.$$

but these are outside the scope of this course. An inelegant but straightforward proof is to examine each component. The x -component of left hand side reads

$$\frac{\partial}{\partial y} \left(\frac{\partial A_y(x, y, z)}{\partial x} - \frac{\partial A_x(x, y, z)}{\partial y} \right) - \frac{\partial}{\partial z} \left(\frac{\partial A_x(x, y, z)}{\partial z} - \frac{\partial A_z(x, y, z)}{\partial x} \right) = -A_x^{(0,0,2)}(x, y, z) - A_x^{(0,2,0)}(x, y, z) + A_z^{(1,0,1)}(x, y, z) + A_y^{(1,1,0)}(x, y, z)$$

The x -component of right hand side is

$$\frac{\partial}{\partial x} \left(\frac{\partial A_x(x, y, z)}{\partial x} + \frac{\partial A_y(x, y, z)}{\partial y} + \frac{\partial A_z(x, y, z)}{\partial z} \right) - \frac{\partial^2 A_x(x, y, z)}{\partial x^2} - \frac{\partial^2 A_x(x, y, z)}{\partial y^2} - \frac{\partial^2 A_x(x, y, z)}{\partial z^2} = -A_x^{(0,0,2)}(x, y, z) - A_x^{(0,2,0)}(x, y, z) + A_z^{(1,0,1)}(x, y, z) + A_y^{(1,1,0)}(x, y, z)$$

These two results are identical


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% == %%  
True
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You can either compute the other two components or argue that x , y , and z are just labels which can be permuted at will (which implies that proving the identity for one component proves it for all components).

Note that $\nabla^2 \mathbf{A}$ consists of a scalar operator (∇^2) applied to a vector (\mathbf{A}) resulting in a vector.

Appendix A

Differential Operators

A.1 Definitions

When parameterizing the position of a point in 3-space, the particular coordinate system which is most useful depends on the symmetry of the physical or geometric system at hand. All coordinate systems can be derived from the Cartesian system by a particular (non-linear) transformation. If $\mathbf{x} = \sum_j \mathbf{e}_j x_j$ is a particular geometric point referred to a rectangular frame of reference, the same point may also be described by coordinates q_i derived from the transformation

$$x_i = x_i(q_j), \quad i, j = 1, 2, 3. \quad (\text{A.1})$$

To obtain the transformation between different coordinate systems, one needs to compute the partial derivatives $\partial x_i / \partial q_j$. The matrix of partial derivatives (the linear map $Dx(q)$) is known as the *Jacobian matrix* (of x):

$$Dx(q) \equiv \begin{pmatrix} \frac{\partial x_1}{\partial q_1} & \frac{\partial x_1}{\partial q_2} & \frac{\partial x_1}{\partial q_3} \\ \frac{\partial x_2}{\partial q_1} & \frac{\partial x_2}{\partial q_2} & \frac{\partial x_2}{\partial q_3} \\ \frac{\partial x_3}{\partial q_1} & \frac{\partial x_3}{\partial q_2} & \frac{\partial x_3}{\partial q_3} \end{pmatrix}. \quad (\text{A.2})$$

A.1.1 Orthogonal coordinates

The coordinates q_j comprise an *orthogonal set* if

$$\sum_{i=1}^3 \frac{\partial x_i}{\partial q_j} \frac{\partial x_i}{\partial q_k} = 0, \quad k \neq j. \quad (\text{A.3})$$

A.1.2 Scale factors

Introducing *dimensional scale factors* $h_i(q_j)$ defined by

$$h_j^2 = \sum_{i=1}^3 \left(\frac{\partial x_i}{\partial q_j} \right)^2, \quad j = 1, 2, 3, \quad (\text{A.4})$$

then, for an orthogonal set of coordinates, one finds that

$$(Dx(q))^T \cdot Dx(q) \equiv \begin{pmatrix} \frac{\partial x_1}{\partial q_1} & \frac{\partial x_2}{\partial q_1} & \frac{\partial x_3}{\partial q_1} \\ \frac{\partial x_1}{\partial q_2} & \frac{\partial x_2}{\partial q_2} & \frac{\partial x_3}{\partial q_2} \\ \frac{\partial x_1}{\partial q_3} & \frac{\partial x_2}{\partial q_3} & \frac{\partial x_3}{\partial q_3} \end{pmatrix} \cdot \begin{pmatrix} \frac{\partial x_1}{\partial q_1} & \frac{\partial x_1}{\partial q_2} & \frac{\partial x_1}{\partial q_3} \\ \frac{\partial x_2}{\partial q_1} & \frac{\partial x_2}{\partial q_2} & \frac{\partial x_2}{\partial q_3} \\ \frac{\partial x_3}{\partial q_1} & \frac{\partial x_3}{\partial q_2} & \frac{\partial x_3}{\partial q_3} \end{pmatrix} = \begin{pmatrix} h_1^2 & 0 & 0 \\ 0 & h_2^2 & 0 \\ 0 & 0 & h_3^2 \end{pmatrix} \quad (\text{A.5})$$

where T denotes the matrix transpose.

A.1.3 Jacobian determinant

Since $|A^T| = |A|$ and $|A \cdot B| = |A| |B|$, where $|A|$ denotes the determinant of the matrix A , we find that the *Jacobian determinant*, $Jx(q)$, for an orthogonal set of coordinates is

$$Jx(q) = |Dx(q)| \equiv \frac{\partial(x_1, x_2, x_3)}{\partial(q_1, q_2, q_3)}(q_1, q_2, q_3) = h_1 h_2 h_3. \quad (\text{A.6})$$

A.1.4 Taylor expansion

The Taylor expansion of $f = \{f_1, f_2, \dots, f_m\} : \mathbb{R}^n \rightarrow \mathbb{R}^m$ can be written as

$$f(y) = f(x) + Df(x) \cdot (y - x) + \frac{1}{2!} D^2 f(x) \cdot (y - x, y - x) + \dots + \frac{1}{k!} D^k f(x) \cdot (y - x, \dots, y - x) + \dots \quad (\text{A.7})$$

where the *Jacobian matrix* of f is

$$Df(x) \equiv \begin{pmatrix} \frac{\partial f_1}{\partial x_1} & \frac{\partial f_1}{\partial x_2} & \dots & \frac{\partial f_1}{\partial x_n} \\ \frac{\partial f_2}{\partial x_1} & \frac{\partial f_2}{\partial x_2} & \dots & \frac{\partial f_2}{\partial x_n} \\ \vdots & \vdots & \ddots & \vdots \\ \frac{\partial f_m}{\partial x_1} & \frac{\partial f_m}{\partial x_2} & \dots & \frac{\partial f_m}{\partial x_n} \end{pmatrix}, \quad (\text{A.8})$$

with $(i, j)^{\text{th}}$ element $\{Df(x)\}_{i,j} = \frac{\partial f_i}{\partial x_j}$ and $D^k f(x) \cdot (y - x, \dots, y - x)$ denotes $D^k f(x)$ as a k -linear map applied to the k -tuple $(y - x, \dots, y - x)$. In coordinates,

$$D^k f(x) (y - x, \dots, y - x) = \sum_{i_1, \dots, i_k=1}^n \left(\frac{\partial^k f}{\partial x_{i_1} \dots \partial x_{i_k}} \right) (y_{i_1} - x_{i_1}) \dots (y_{i_k} - x_{i_k}). \quad (\text{A.9})$$

A.1.5 Differential elements

The differential elements are

$$\begin{aligned} \text{Line :} & \quad ds_k = h_k dq_k \\ \text{Area :} & \quad dS_{j,k} = dx_j dx_k \rightarrow h_j h_k dq_j dq_k \\ \text{Volume :} & \quad dV = dx dy dz \rightarrow Jx(q) dq_1 dq_2 dq_3 = h_1 h_2 h_3 dq_1 dq_2 dq_3 \end{aligned} \quad (\text{A.10})$$

A.1.6 Differential operators

The fundamental vector operators can be shown to be

$$\text{Gradient : } \nabla f = \frac{1}{h_1} \frac{\partial f}{\partial q_1} \mathbf{e}_1 + \frac{1}{h_2} \frac{\partial f}{\partial q_2} \mathbf{e}_2 + \frac{1}{h_3} \frac{\partial f}{\partial q_3} \mathbf{e}_3 \quad (\text{A.11})$$

$$\begin{aligned} \text{Divergence : } \nabla \cdot \mathbf{F} = & \\ \frac{1}{Jx(q)} \left(\frac{\partial}{\partial q_1} \left(\frac{Jx(q)}{h_1} F_1 \right) + \frac{\partial}{\partial q_2} \left(\frac{Jx(q)}{h_2} F_2 \right) + \frac{\partial}{\partial q_3} \left(\frac{Jx(q)}{h_3} F_3 \right) \right) & \quad (\text{A.12}) \end{aligned}$$

$$\text{Laplacian: } \nabla^2 f = \frac{1}{Jx(q)} \left(\frac{\partial}{\partial q_1} \left(\frac{Jx(q)}{h_1^2} \frac{\partial f}{\partial q_1} \right) + \frac{\partial}{\partial q_2} \left(\frac{Jx(q)}{h_2^2} \frac{\partial f}{\partial q_2} \right) + \frac{\partial}{\partial q_3} \left(\frac{Jx(q)}{h_3^2} \frac{\partial f}{\partial q_3} \right) \right) \quad (\text{A.13})$$

$$\text{Curl: } \nabla \wedge \mathbf{F} = \frac{1}{Jx(q)} \begin{vmatrix} h_1 \mathbf{e}_1 & h_2 \mathbf{e}_2 & h_3 \mathbf{e}_3 \\ \frac{\partial}{\partial q_1} & \frac{\partial}{\partial q_2} & \frac{\partial}{\partial q_3} \\ h_1 F_1 & h_2 F_2 & h_3 F_3 \end{vmatrix} \quad (\text{A.14})$$

A.2 Examples

A.2.1 Spherical Polar Coordinates

A.2.1.1 Definition

$$x = r \cos(\phi) \sin(\theta), \quad y = r \sin(\phi) \sin(\theta), \quad z = r \cos(\theta) \quad (\text{A.15})$$

Note that care must be taken when inverting these relations. For example, $\phi \neq \tan^{-1}(\frac{y}{x})$ in general — though you will see this statement appearing regularly in textbooks.

A.2.1.2 Jacobian Matrix

From the definition, we find

$$\mathcal{D} = \begin{pmatrix} \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial r} & \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial \theta} & \frac{\partial(r \cos(\phi) \sin(\theta))}{\partial \phi} \\ \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial r} & \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial \theta} & \frac{\partial(r \sin(\phi) \sin(\theta))}{\partial \phi} \\ \frac{\partial(r \cos(\theta))}{\partial r} & \frac{\partial(r \cos(\theta))}{\partial \theta} & \frac{\partial(r \cos(\theta))}{\partial \phi} \end{pmatrix} \\ \begin{pmatrix} \cos(\phi) \sin(\theta) & r \cos(\theta) \cos(\phi) & -r \sin(\theta) \sin(\phi) \\ \sin(\theta) \sin(\phi) & r \cos(\theta) \sin(\phi) & r \cos(\phi) \sin(\theta) \\ \cos(\theta) & -r \sin(\theta) & 0 \end{pmatrix}$$

This matrix can also be generated directly by taking the outer product of the partial derivative (\mathcal{D}) with the coordinate vectors:

$$\text{Outer}[\mathcal{D}, \{r \cos(\phi) \sin(\theta), r \sin(\phi) \sin(\theta), r \cos(\theta)\}, \{r, \theta, \phi\}] \\ \begin{pmatrix} \cos(\phi) \sin(\theta) & r \cos(\theta) \cos(\phi) & -r \sin(\theta) \sin(\phi) \\ \sin(\theta) \sin(\phi) & r \cos(\theta) \sin(\phi) & r \cos(\phi) \sin(\theta) \\ \cos(\theta) & -r \sin(\theta) & 0 \end{pmatrix}$$

A.2.1.3 Jacobian Determinant

The Jacobian determinant is needed when computing the volume element:

$$\mathcal{J} = |\mathcal{D}| \text{ // Simplify} \\ r^2 \sin(\theta)$$

A.2.1.4 Scale Factors

Computing

$$\mathcal{D}^T \cdot \mathcal{D} \text{ // Simplify}$$

$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & r^2 & 0 \\ 0 & 0 & r^2 \sin^2(\theta) \end{pmatrix}$$

we see that spherical polar coordinates are an orthogonal coordinate system. The diagonal entries,

$$\text{Transpose}(\%, \{1, 1\})$$

$$\{1, r^2, r^2 \sin^2(\theta)\}$$

lead to the scale factors:

$$\{h_r, h_\theta, h_\phi\} = \sqrt{\%} \text{ // PowerExpand}$$

$$\{1, r, r \sin(\theta)\}$$

A.2.1.5 Differential Operators

From the definitions, we find that

$$\text{Gradient: } \nabla f = \frac{\partial f}{\partial r} \mathbf{e}_r + \frac{1}{r} \frac{\partial f}{\partial \theta} \mathbf{e}_\theta + \frac{1}{r \sin(\theta)} \frac{\partial f}{\partial \phi} \mathbf{e}_\phi \quad (\text{A.16})$$

$$\text{Divergence: } \nabla \cdot \mathbf{F} = \frac{1}{r^2} \frac{\partial}{\partial r} (r^2 F_r) + \frac{1}{r \sin(\theta)} \frac{\partial}{\partial \theta} (\sin(\theta) F_\theta) + \frac{1}{r \sin(\theta)} \frac{\partial F_\phi}{\partial \phi} \quad (\text{A.17})$$

$$\text{Laplacian: } \nabla^2 f = \frac{1}{r^2} \frac{\partial}{\partial r} \left(r^2 \frac{\partial f}{\partial r} \right) + \frac{1}{r^2} \left(\frac{1}{\sin(\theta)} \frac{\partial}{\partial \theta} \left(\sin(\theta) \frac{\partial f}{\partial \theta} \right) + \frac{1}{\sin(\theta)^2} \frac{\partial^2 f}{\partial \phi^2} \right) \quad (\text{A.18})$$

$$\text{Curl: } \nabla \wedge \mathbf{F} =$$

$$\frac{1}{r^2 \sin(\theta)} \begin{vmatrix} \mathbf{e}_r & r \mathbf{e}_\theta & r \sin(\theta) \mathbf{e}_\phi \\ \partial_r & \partial_\theta & \partial_\phi \\ F_r & r F_\theta & r \sin(\theta) F_\phi \end{vmatrix} = \frac{1}{r \sin(\theta)} \left(\frac{\partial(\sin(\theta) F_\phi)}{\partial \theta} - \frac{\partial F_\theta}{\partial \phi} \right) \mathbf{e}_r +$$

$$\frac{1}{r \sin(\theta)} \left(\frac{\partial F_r}{\partial \phi} - \sin(\theta) \frac{\partial(r F_\phi)}{\partial r} \right) \mathbf{e}_\theta + \frac{1}{r} \left(\frac{\partial(r F_\theta)}{\partial r} - \frac{\partial F_r}{\partial \theta} \right) \mathbf{e}_\phi \quad (\text{A.19})$$

A.2.2 Cylindrical Coordinates

A.2.2.1 Definition

$$x = \rho \cos(\phi), \quad y = \rho \sin(\phi), \quad z = z \quad (\text{A.20})$$

Note that care must be taken when inverting these relations. For example, $\phi \neq \tan^{-1}(\frac{y}{x})$ in general — though you will see this statement appearing regularly in textbooks. However, using half-angle formulæ, you can show that $\phi = 2 \tan^{-1}(\frac{y}{x+\rho})$ is correct. Alternatively, you can use $\phi = \tan^{-1}(y/x)$, a special form of \tan^{-1} which gives the arc tangent of y/x , taking into account which quadrant the point (x, y) is in.

$$\frac{y}{x + \rho} /. \{x \rightarrow \rho \cos(\phi), y \rightarrow \rho \sin(\phi)\} // \text{Simplify}$$

$$\tan\left(\frac{\phi}{2}\right)$$

A.2.2.2 Jacobian Matrix

We generate the Jacobian matrix directly by taking the outer product of the partial derivative (\mathcal{D}) of the coordinate vectors:

$$\mathcal{D} = \text{Outer}[\mathcal{D}, \{\rho \cos(\phi), \rho \sin(\phi), z\}, \{\rho, \phi, z\}]$$

$$\begin{pmatrix} \cos(\phi) & -\rho \sin(\phi) & 0 \\ \sin(\phi) & \rho \cos(\phi) & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

A.2.2.3 Jacobian Determinant

The Jacobian determinant is needed when computing the volume element:

$$\mathcal{J} = |\mathcal{D}| // \text{Simplify}$$

$$\rho$$

A.2.2.4 Scale Factors

Computing

$$\mathcal{D}^T \cdot \mathcal{D} // \text{Simplify}$$

$$\begin{pmatrix} 1 & 0 & 0 \\ 0 & \rho^2 & 0 \\ 0 & 0 & 1 \end{pmatrix}$$

we see that spherical polar coordinates are an orthogonal coordinate system. The diagonal entries,

$$\text{Transpose}(\%, \{1, 1\})$$

$$\{1, \rho^2, 1\}$$

lead to the scale factors:

$$\{h_\rho, h_\phi, h_z\} = \sqrt{\%} // \text{PowerExpand}$$

$$\{1, \rho, 1\}$$

A.2.2.5 Differential Operators

From the definitions, we find that

$$\text{Gradient} : \nabla f = \frac{\partial f}{\partial \rho} \mathbf{e}_\rho + \frac{1}{\rho} \frac{\partial f}{\partial \phi} \mathbf{e}_\phi + \frac{\partial f}{\partial z} \mathbf{e}_z \quad (\text{A.21})$$

$$\text{Divergence} : \nabla \cdot \mathbf{F} = \frac{1}{\rho} \frac{\partial(\rho F_\rho)}{\partial \rho} + \frac{1}{\rho} \frac{\partial F_\phi}{\partial \phi} + \frac{\partial F_z}{\partial z} \quad (\text{A.22})$$

$$\text{Laplacian : } \nabla^2 f = \frac{1}{\rho} \left(\frac{\partial}{\partial \rho} \left(\rho \frac{\partial f}{\partial \rho} \right) \right) + \frac{1}{\rho^2} \frac{\partial^2 f}{\partial \phi^2} + \frac{\partial^2 f}{\partial z^2} \quad (\text{A.23})$$

$$\text{Curl : } \nabla \wedge \mathbf{F} = \frac{1}{\rho} \begin{vmatrix} \mathbf{e}_\rho & \rho \mathbf{e}_\phi & \mathbf{e}_z \\ \partial_\rho & \partial_\phi & \partial_z \\ F_\rho & \rho F_\phi & F_z \end{vmatrix} = \quad (\text{A.24})$$

$$\left(\frac{1}{\rho} \frac{\partial F_z}{\partial \phi} - \frac{\partial F_\phi}{\partial z} \right) \mathbf{e}_\rho + \left(\frac{\partial F_\rho}{\partial z} - \frac{\partial F_z}{\partial \rho} \right) \mathbf{e}_\phi + \frac{1}{\rho} \left(\frac{\partial(\rho F_\phi)}{\partial \rho} - \frac{\partial F_\rho}{\partial \phi} \right) \mathbf{e}_z$$

Appendix B

Dirac's Delta Function $\delta(x)$

B.1 Examples

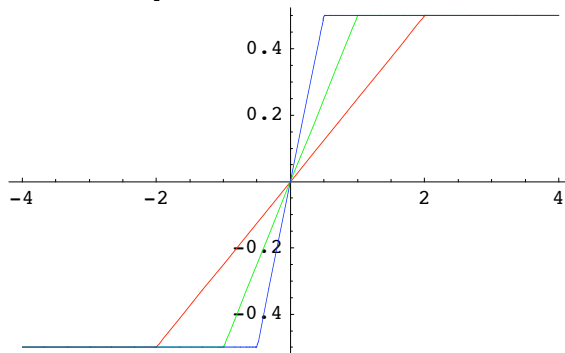
Consider the (piecewise continuous) function

$$\theta_{\epsilon}[x] := -\frac{1}{2} /; x < -\epsilon$$

$$\theta_{\epsilon}[x] := \frac{x}{2\epsilon} /; -\epsilon \leq x \leq \epsilon$$

$$\theta_{\epsilon}[x] := \frac{1}{2} /; x > \epsilon$$

$$\text{Plot}\left\{\theta_2(x), \theta_1(x), \theta_{\frac{1}{2}}(x)\right\}, \{x, -4, 4\}, \text{PlotStyle} \rightarrow \{\text{Hue}(0), \text{Hue}\left(\frac{1}{3}\right), \text{Hue}\left(\frac{2}{3}\right)\};$$



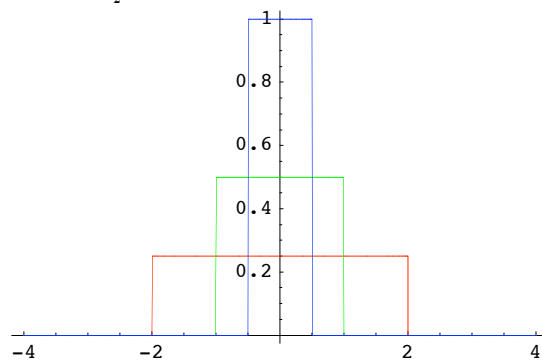
The derivative is

$$\delta_{\epsilon}[x] := 0 /; x < -\epsilon$$

$$\delta_{\epsilon}[x] := \frac{1}{2\epsilon} /; -\epsilon \leq x \leq \epsilon$$

$$\delta_{\epsilon}[x] := 0 /; x > \epsilon$$

$$\text{Plot}\left\{\delta_2(x), \delta_1(x), \delta_{\frac{1}{2}}(x)\right\}, \{x, -4, 4\}, \text{PlotStyle} \rightarrow \{\text{Hue}(0), \text{Hue}\left(\frac{1}{3}\right), \text{Hue}\left(\frac{2}{3}\right)\};$$



The area under each curve is *constant*:

$$\mathbf{NIntegrate}[\delta_1(x), \{x, -5, -1, 1, 5\}]$$

1.

$$\mathbf{NIntegrate}[\delta_{1/2}(x), \{x, -5, -1/2, 1/2, 5\}]$$

1.

In other words

$$\int_{-\infty}^{\infty} \delta_{\epsilon}(x) dx = 1, \quad \forall \epsilon \in \mathbb{R}^+.$$

Alternatively, we can obtain this from

$$\int_{-\infty}^{\infty} \delta_{\epsilon}(x) dx = \int_{-\infty}^{\infty} \theta_{\epsilon}'(x) dx = \theta_{\epsilon}(x)|_{-\infty}^{\infty} = \left(\frac{1}{2} - \left(-\frac{1}{2}\right)\right) = 1, \quad \forall \epsilon \in \mathbb{R}^+.$$

Consider the integral

$$\int_{-\infty}^{\infty} f(x) \delta_{\epsilon}(x) dx,$$

where $f(x)$ is an arbitrary function which goes to 0 "sufficiently fast" as $x \rightarrow \pm\infty$. For ϵ "sufficiently small", we can write

$$\int_{-\infty}^{\infty} f(x) \delta_{\epsilon}(x) dx = \frac{1}{2\epsilon} \int_{-\epsilon}^{\epsilon} f(x) dx \approx \frac{1}{2\epsilon} (2\epsilon f(0)) = f(0),$$

using the Mean Value Theorem. As $\epsilon \rightarrow 0$, $\delta_{\epsilon}(x) \rightarrow \delta(x)$, and $\delta(x)$ has the interesting properties that

$$\int_{-\infty}^{\infty} \delta(x) dx = 1, \quad \int_{-\infty}^{\infty} f(x) \delta(x) dx = f(0)$$

Show, using integration by parts, that

$$\int_{-\infty}^{\infty} f(x) \delta'(x) dx = -f'(0)$$

B.2 Definition

The symbol $\delta(x)$ is not a function in the usual mathematical sense. A function in one dimension is a mapping between ordered pairs $x \rightarrow y = f(x)$. In the case of the symbol $\delta(x)$, any such mapping carries every point x on the real axis, save one, into the number zero. This is hardly a well behaved function. Nevertheless, it can be treated symbolically as though it shared most properties of ordinary smooth

functions. I will often treat it as an ordinary function. Our purpose here is to outline this highly useful notation, and not to give a mathematical justification for this use.

It is possible to view $\delta(x)$ as representing the symbolic limit of a sequence of suitably defined functions. Imagine, for example, a sequence based on the parameter ϵ defined by any of the following three functions:

Code

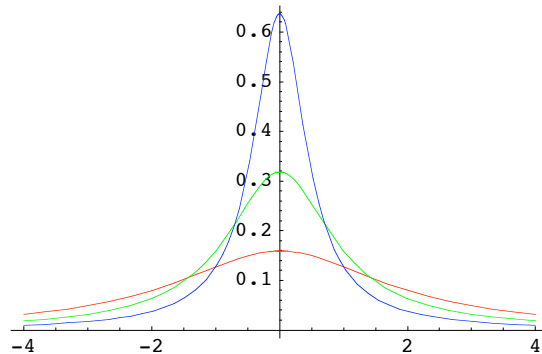
Clear[Subscript]

SetOptions[Integrate, GenerateConditions → False];

$$\delta_\epsilon(x) = \frac{\epsilon}{\pi} \frac{1}{x^2 + \epsilon^2}$$

$$\delta_\epsilon[x_] := \frac{\epsilon}{\pi} \frac{1}{x^2 + \epsilon^2}$$

Plot $\left(\{\delta_2(x), \delta_1(x), \delta_{\frac{1}{2}}(x)\}, \{x, -4, 4\}, \text{PlotStyle} \rightarrow \{\text{Hue}(0), \text{Hue}\left(\frac{1}{3}\right), \text{Hue}\left(\frac{2}{3}\right)\}\right);$



For any ϵ we find that

$$\int_{-\infty}^{\infty} \delta_\epsilon(x) dx // \text{PowerExpand}$$

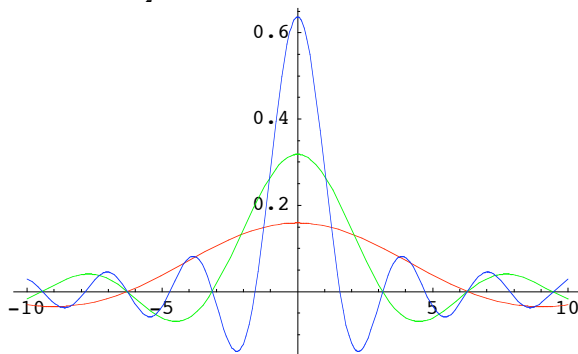
For the function $\frac{1}{1+x^2}$

$$\lim_{\epsilon \rightarrow 0} \int_{-\infty}^{\infty} \frac{1}{1+x^2} \delta_\epsilon(x) dx // \text{PowerExpand}$$

$$\delta_\epsilon(x) = \frac{1}{\pi} \frac{\sin\left(\frac{x}{\epsilon}\right)}{x}$$

$$\delta_\epsilon[x_] := \frac{1}{\pi} \frac{\sin\left(\frac{x}{\epsilon}\right)}{x}$$

Plot $\left(\{\delta_2(x), \delta_1(x), \delta_{\frac{1}{2}}(x)\}, \{x, -10, 10\}, \text{PlotRange} \rightarrow \text{All}, \text{PlotStyle} \rightarrow \{\text{Hue}(0), \text{Hue}\left(\frac{1}{3}\right), \text{Hue}\left(\frac{2}{3}\right)\}\right);$

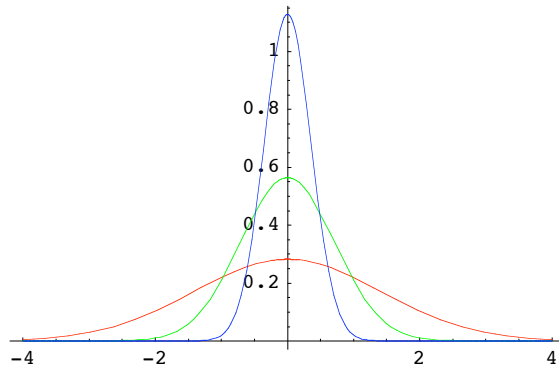


$$\int_{-\infty}^{\infty} \delta_{\epsilon}(x) dx / \text{sgn}(\epsilon) \rightarrow 1$$

$$\delta_{\epsilon}(x) = \frac{1}{\sqrt{\pi}} \frac{1}{\epsilon} e^{-\frac{x^2}{\epsilon^2}}$$

$$\delta_{\epsilon}[x_] := \frac{1}{\sqrt{\pi}} \frac{1}{\epsilon} e^{-\frac{x^2}{\epsilon^2}}$$

`Plot`{ $\{\delta_2(x), \delta_1(x), \delta_{\frac{1}{2}}(x)\}$, { $x, -4, 4$ }, `PlotStyle` \rightarrow {`Hue`(0), `Hue`($\frac{1}{3}$), `Hue`($\frac{2}{3}$)}, `PlotRange` \rightarrow `All`};



$$\int_{-\infty}^{\infty} \delta_{\epsilon}(x) dx // \text{PowerExpand}$$

B.3 Sequence

Each of these functions, for any $\epsilon \ll 1$, have the properties (1) sharply peaked at $x = 0$; and (2) area under curve is unity independent of ϵ . In short, if one constructs a convergent sequence of ϵ 's, then the quantity $\delta(x) = \lim_{\epsilon \rightarrow 0} \delta_{\epsilon}(x)$.

for each of the above functions $\delta_{\epsilon}(x)$, has the desired properties of the delta "function". By this I mean the following: whenever a delta "function" appears multiplying a smooth function under an integral sign, you should imagine that it is replaced by $\delta_{\epsilon}(x)$ and the integral evaluated. Then, after integration, the limit of the sequence of the results of integration is taken. This process gives meaning to the delta function. With this idea in mind, I can treat the delta function as if it were itself a smooth function and even write, for example,

$$\int_{-\infty}^{\infty} \delta'(x) f(x) dx = - \int_{-\infty}^{\infty} \delta(x) f'(x) dx = -f'(0)$$

where I have integrated by parts. Other useful results, whose justification are based in such arguments, are these:

$$\begin{aligned} \delta(ax) &= \frac{\delta(x)}{|a|} \\ x \delta(x) &= 0 \\ \delta(x^2 - a^2) &= \frac{1}{2|a|} (\delta(x - a) + \delta(x + a)) \\ \delta((x - a)(x - b)) &= \frac{1}{|b - a|} (\delta(x - a) + \delta(x - b)) \end{aligned}$$

B.4 Three Dimensions

The idea of the delta function is readily extended to spaces of higher dimension. Let a point in 3-dimensions be $\mathbf{r} = (x, y, z)$. I define the delta function in this space by

$$\begin{aligned}\delta(\mathbf{r} - \mathbf{r}') &= \delta(x - x') \delta(y - y') \delta(z - z') \\ \int \int \int \delta(\mathbf{r} - \mathbf{r}') dx dy dz &= 1, \\ \int \int \int \delta(\mathbf{r} - \mathbf{r}') f(\mathbf{r}) dx dy dz &= f(\mathbf{r}')\end{aligned}$$

An example of this is

$$\int \int \int \delta(x - a) e^{-|x-b|} dx dy dz = e^{-|b-a|}.$$

As a further example, consider

$$h(\mathbf{r}) = \nabla^2 \left(\frac{1}{r} \right) = \nabla \cdot \nabla \left(\frac{1}{r} \right) = -\nabla \cdot \frac{\mathbf{r}}{r^3} = -\left(\frac{1}{r^3} \nabla \cdot \mathbf{r} - 3 \mathbf{r} \cdot \left(\frac{\mathbf{r}}{r^5} \right) \right) = -\left(\frac{3}{r^3} - \frac{3}{r^3} \right) = 0$$

if $r \neq 0$.

Clearly, $h(\mathbf{r})$ is singular at $r = 0$. To investigate its behavior near the singularity, integrate over a small spherical volume V centered at the origin. The divergence theorem

$$\int_V \nabla \cdot \mathbf{A} d\mathbf{r} = \int_{S=\partial V} \mathbf{A} \cdot d\mathbf{S},$$

provides what we seek:

$$\begin{aligned}\int_V h(\mathbf{r}) d\mathbf{r} &= \int_V \nabla^2 \left(\frac{1}{r} \right) d\mathbf{r} \\ &= \int_V \nabla \cdot \nabla \left(\frac{1}{r} \right) d\mathbf{r} \\ &= \int_{S=\partial V} \nabla \left(\frac{1}{r} \right) \cdot d\mathbf{S} \\ &= \int_{S=\partial V} \nabla \left(\frac{1}{r} \right) \cdot \hat{\mathbf{r}} dS \\ &= -\int_{S=\partial V} \frac{r}{r^3} \cdot \hat{\mathbf{r}} r^2 d\Omega \\ &= -\int_{S=\partial V} d\Omega \\ &= -4\pi.\end{aligned}$$

Conclusion: $h(\mathbf{r})$ is zero everywhere except at a single point, namely $r = 0$. There it is infinite, but in such a way that its volume integral over the singularity is -4π . Therefore, we have the identity

$$-\nabla^2 \left(\frac{1}{r} \right) = \nabla \cdot \frac{\hat{\mathbf{r}}}{r^2} = 4\pi \delta(\mathbf{r}).$$