

6 Tensors

A famously annoying definition of a tensor is:

A tensor is something whose components transform like a tensor

This becomes even more annoying when you appreciate that this is, in fact, one of the better definitions of a tensor. The purpose of this section is to explain why this definition is not as dumb as it sounds and to give some insight into what it means to be a tensor.

Very roughly speaking, tensors are generalisations of objects like vectors and matrices. In index notation, a vector has a single index while a matrix has two indices. A tensor is an object with any number of indices, something like $T_{ij\dots k}$.

However, this simplistic description hides the most important property of a tensor. Vectors, matrices and, more generally, tensors are more than just a list of numbers. Instead, those numbers should be thought of as a useful way of characterising the underlying object and, because of this, inherit some properties of that underlying object. As we will see, the key property is how the list of numbers transform under a change of basis.

We will start by explaining this in more detail, firstly with vectors and then building up to the definition of a tensor. Initially we will keep the discussion restricted to some (admittedly rather dry) mathematical formalism. Then, in Section 6.2 we will describe some physical examples.

6.1 What it Takes to Make a Tensor

Not any list of n numbers constitutes a vector in \mathbb{R}^n . Or, said more precisely, not any list of n numbers constitutes the components of a vector in \mathbb{R}^n . For example, if you write down the heights of the first three people you met this morning, that doesn't make a vector in \mathbb{R}^3 . Instead, a vector comes with certain responsibilities. In particular, the components describe an underlying object which should be independent of the choice of basis. As we now explain, that means that the components should transform in the right way under rotations.

We consider a point $\mathbf{x} \in \mathbb{R}^n$. If we wish to attach some coordinates to this point, we first need to introduce a set of basis vectors $\{\mathbf{e}_i\}$ with $i = 1, \dots, n$. We will take these to be orthonormal, meaning that $\mathbf{e}_i \cdot \mathbf{e}_j = \delta_{ij}$. Any vector can then be expressed as

$$\mathbf{x} = x_i \mathbf{e}_i \tag{6.1}$$

Usually we conflate the components $x_i = (x_1, \dots, x_n)$ with the “vector”. But, for our purposes, we should remember that these are just a useful way of representing the more abstract object \mathbf{x} . In particular, we’re entirely at liberty to take a different set of basis vectors,

$$\mathbf{e}'_i = R_{ij}\mathbf{e}_j$$

If we ask that \mathbf{e}'_i are also orthonormal, so $\mathbf{e}'_i \cdot \mathbf{e}'_j = \delta_{ij}$, then we have

$$\mathbf{e}'_i \cdot \mathbf{e}'_j = R_{ik}R_{jl}\mathbf{e}_k \cdot \mathbf{e}_l = R_{ik}R_{jk} = \delta_{ij}$$

or, in matrix notation,

$$RR^T = \mathbf{1}$$

Matrices of this kind are said to be *orthogonal*. We write $R \in O(n)$. Taking the determinant, we have $\det R = \pm 1$. Those matrices with $\det R = +1$ correspond to rotations and are said to be *special orthogonal*. We write $R \in SO(n)$. In \mathbb{R}^3 , a rotation $R \in SO(3)$ takes a right-handed orthonormal basis into another right-handed orthonormal basis. Those matrices with $\det R = -1$ correspond to a rotation together with a reflection and take a right-handed basis to a left-handed basis.

Under a change of basis, the vector \mathbf{x} itself doesn’t change. But its components do. We have

$$\mathbf{x} = x_i\mathbf{e}_i = x'_i\mathbf{e}'_i = x'_iR_{ij}\mathbf{e}_j$$

So the components transform under the same rotation matrix R ,

$$x_j = R_{ij}x'_i \quad \Rightarrow \quad x'_i = R_{ij}x_j \quad (6.2)$$

A *tensor* T is a generalisation of these ideas to an object with more indices. Just as the vector \mathbf{x} has an identity independent of any choice of basis, so too does the tensor T . But when measured with respect to a chosen basis $\{\mathbf{e}_i\}$, a *tensor of rank p* has components $T_{i_1 \dots i_p}$. When we change the basis using (6.1), the tensor transforms as

$$T'_{i_1 \dots i_p} = R_{i_1 j_1} \dots R_{i_p j_p} T_{j_1 \dots j_p} \quad (6.3)$$

This is known as the *tensor transformation rule*. A tensor of rank p is sometimes referred to simply as a p -tensor.

The simplest examples of tensors are very familiar. A tensor of rank 0 is just a number, or scalar, T . Here there's no requirement because a number doesn't change if you do a rotation: $T' = T$. So any single number can be said to be a tensor, although it isn't a particularly helpful designation.

A tensor of rank 1 is a vector. Here, however, it's important that the components of the vector transform as $T'_j = R_{ij}T_j$. If they don't transform in this way, then you don't have a tensor on your hands. You just have a bunch of numbers.

A tensor of rank 2 is a matrix that transforms as $T'_{ij} = R_{ik}R_{jl}T_{kl}$. Again, the transformation property is key. Just because you have an array of numbers A_{ij} , arranged in an $n \times n$ grid, doesn't mean that you have a 2-tensor. You have to check the transformation property holds. Otherwise, as with a vector, the array of numbers isn't a tensor; it's just a bunch of numbers.

What's a Tensor and What's Not?

It's worth elaborating on the definition of a tensor. For example, suppose that someone hands you a matrix, say

$$T_{ij} = \begin{pmatrix} 3 & 8 & 0 \\ 5 & -4 & 3 \\ 1 & 1 & 3 \end{pmatrix}$$

and asks you: "is this a tensor?". It's natural to answer yes. After all, it's written as T_{ij} which is the name we've given to a tensor. And it looks for all the world like a matrix. So is it a tensor? The answer is: we don't know. We haven't been given enough information. As we've stressed several times, a tensor isn't just a bunch of numbers arranged in some pattern. This sometimes goes by the name of an *array* of numbers. Instead, we only know that a given array of numbers is a tensor if it transforms as (6.3). That means that we need to firstly know what basis the array of numbers above has been measured in. And then we need to know what the array looks like when measured in other bases. Only then do we have enough information to say whether this is a tensor or not. It's a tensor only if transforms as (6.3): this transformation law is the definition of a tensor.

Here's another example. In a given basis, the position of a point is given by x_i . We write this as the components of a vector

$$x_i = \begin{pmatrix} x \\ y \\ z \end{pmatrix}$$

This is a tensor. Indeed, our starting point is that the components of this simple vector transforms in the tensorial way (6.2). This is just the statement that the components of this vector transform in the familiar way under rotation.

Suppose that you now square each of these elements and decide to write them as a column vector. We'll give it a fancy name Λ_i , complete with that hanging i index,

$$\Lambda_i = \begin{pmatrix} x^2 \\ y^2 \\ z^2 \end{pmatrix}$$

That i index makes this look for all the world like it's a tensor. But it's not. We know that after a rotation, $x_i \rightarrow x'_i = R_{ij}x_j$. This means that if we do a rotation and then measure the components of the array Λ'_i we get

$$\Lambda'_i = \begin{pmatrix} (R_{11}x + R_{12}y + R_{13}z)^2 \\ (R_{21}x + R_{22}y + R_{23}z)^2 \\ (R_{31}x + R_{32}y + R_{33}z)^2 \end{pmatrix}$$

But that's most definitely not how a tensor transforms! It's not the rule (6.3) that we wanted. The upshot is that Λ_i is not a tensor and it was a little bit naughty to write it as Λ_i because it suggests that it has some property that it doesn't.

Relatedly, this explains something that you may have wondered about in school. Suppose that you're given two vectors. You know that you can take an inner product to get a scalar, or you can take the cross-product to get another vector. But what stops you from doing something much simpler, just multiplying the component of one vector with the corresponding component of another vector to get a third vector. It seems like such an obvious thing to do. But it's a bad thing to do, precisely because the thing you end up with is not a tensor. It does not transform in the way (6.2), which is how components of a vector should transform.

There is a similar story for matrices. If you have two matrices, then there’s a ridiculously complicated way to multiply them, multiplying rows with columns. Why don’t we just do something much simpler and multiply entries together component by component? You’ve probably guessed the answer by now. If we started with genuine matrices, meaning that they transform (6.3), then the object that you get if you do proper matrix multiplication will also transform as (6.3), but the simpler, stupid way to multiplying will not.

Why are we making such a big deal about this? What is so special about things that transform nicely as (6.3) under rotations? Well, there are several answers to this, depending on taste. At the most basic level, if you’re a physicist, then you might genuinely want to know how something looks in different, rotated frames of reference.

Moreover, once you realise that there’s a preferred way for things to transform — the tensor way (6.3) — this brings some extra power to the calculations, a little like dimensional analysis. Suppose that you have an equation of the form “left-hand side” = “right-hand side”. If the thing on the left is a tensor then the thing on the right better also be a tensor. And sometimes there’s not many tensors available, which limits your options for what the thing on the right can actually be. We’ll see an example of this in Section 6.1.3 when we’ll use tensors to make some scary looking integrals a little more palatable.

The discussion above is very much from a physics perspective. But what about a pure maths perspective? This gives a more formal, but arguably cleaner, definition of a tensor. We’ll explain this imminently in Section 6.1.1.

We’ll meet a number of tensors as we proceed. But there is a one that is special: this is the rank 2 tensor δ_{ij} or, equivalently, the unit matrix. Importantly, it has the same 0 and 1 entries in any basis because, under the transformation (6.3), it becomes

$$\delta'_{ij} = R_{ik}R_{jl}\delta_{kl} = \delta_{ij}$$

We will devote Section 6.1.3 to “invariant tensors” which, like δ_{ij} , take the same form in any basis.

6.1.1 Tensors as Maps

There is something a little strange about the definition of a tensor given above. We first pick a set of coordinates, and the transformation law (6.3) then requires that the tensor transforms nicely so that, ultimately, nothing depends on these coordinates. But, if that’s the case, surely there should be a definition of a tensor that doesn’t rely on coordinates at all!

There is. A tensor T of rank p is a multi-linear map that takes p vectors, $\mathbf{a}, \mathbf{b}, \dots, \mathbf{c}$ and spits out a number in \mathbb{R} ,

$$T(\mathbf{a}, \mathbf{b}, \dots, \mathbf{c}) = T_{i_1 i_2 \dots i_p} a_{i_1} b_{i_2} \dots c_{i_p} \quad (6.4)$$

Here “multi-linear” means that T is linear in each of the entries $\mathbf{a}, \mathbf{b}, \dots, \mathbf{c}$ individually. By evaluating T on all possible vectors $\mathbf{a}, \mathbf{b}, \dots, \mathbf{c}$, we get the components $T_{i_1 i_2 \dots i_p}$. The transformation rule (6.3) is simply the statement that the map T is independent of the choice of basis, and we can equally well write

$$\begin{aligned} T(\mathbf{a}, \mathbf{b}, \dots, \mathbf{c}) &= T'_{i_1 i_2 \dots i_p} a'_{i_1} b'_{i_2} \dots c'_{i_p} \\ &= (R_{i_1 j_1} R_{i_2 j_2} \dots R_{i_p j_p} T_{j_1 j_2 \dots j_p})(R_{i_1 k_1} a_{k_1})(R_{i_2 k_2} b_{k_2}) \dots (R_{i_p k_p} c_{k_p}) \\ &= T_{j_1 j_2 \dots j_p} a_{j_1} b_{j_2} \dots c_{j_p} \end{aligned}$$

which follows because $R^T R = \mathbf{1}$ or, in components, $R_{ij} R_{ik} = \delta_{jk}$. The key is that this formula takes the same form in any basis.

Tensors as Maps Between Vectors

Rather than thinking of a tensor as a map from many vectors to \mathbb{R} , you can equivalently think of it as a map from some lower-rank tensor to another. For example, in (6.4), if you don't fill in the first entry, then a rank p tensor can equally well be viewed as taking $(p - 1)$ vectors and spitting out a single vector

$$a_i = T_{i j_1 \dots j_{p-1}} b_{j_1} \dots c_{j_{p-1}}$$

This is the way that tensors typically arise in physics or applied mathematics, where the most common example is simply a rank 2 tensor, defined as a map from one vector to another

$$\mathbf{u} = T\mathbf{v} \quad \Rightarrow \quad u_i = T_{ij} v_j$$

Until now, we've simply called T a matrix but for the equation $\mathbf{u} = T\mathbf{v}$ to make sense, T must transform as a tensor (6.3). This is inherited from the transformation rules of the vectors, $u'_i = R_{ij} u_j$ and $v'_i = R_{ij} v_j$, giving

$$u'_i = T'_{ij} v'_j \quad \text{with} \quad T'_{ij} = R_{ik} R_{jl} T_{kl}$$

Written as a matrix equation, this is $T' = R T R^T$.

6.1.2 Tensor Operations

Given a bunch of tensors, there are some manipulations that leave you with another tensor. Here we describe these operations.

- We can *add* and *subtract* tensors of the same rank, so if S and T are both tensors of rank p then so too is $S + T$. We can also multiply a tensor by a constant α and it remains a tensor.
- If S is a tensor of rank p and T a tensor of rank q , then the *tensor product* $S \otimes T$ is a tensor of rank $p + q$, defined by

$$(S \otimes T)_{i_1 \dots i_p j_1 \dots j_q} = S_{i_1 \dots i_p} T_{j_1 \dots j_q}$$

You can check that the components of $(S \otimes T)$ do indeed satisfy the transformation rule (6.3). In particular, if we have p different vectors \mathbf{a} , \mathbf{b} , \dots , \mathbf{c} then we can construct a tensor

$$T = \mathbf{a} \otimes \mathbf{b} \otimes \dots \otimes \mathbf{c} \quad \text{with} \quad T_{i_1 \dots i_p} = a_{i_1} b_{i_2} \dots c_{i_p}$$

- Given a tensor T of rank p , we can construct a new tensor S of rank $(p - 2)$ by *contracting* on two indices using δ_{ij} ,

$$S_{k_1 \dots k_{p-2}} = \delta_{ij} T_{ijk_1 \dots k_{p-2}}$$

For a rank 2 tensor, the contraction is what we call the trace, $\text{Tr } T = T_{ii}$. It's a valid tensor operation because the end result is a scalar that does not transform under rotations

$$T'_{ii} = R_{ij} R_{ik} T_{jk} = \delta_{jk} T_{jk} = T_{jj}$$

The same derivation shows that higher rank tensors can also be contracted, with the additional indices unaffected by the contraction.

Combining a contraction with a tensor product gives a way to contract two different tensors together. For example, given a p -tensor P and q -tensor Q , we can form a $p + q - 2$ tensor by contracting, say, the first index on each to get $P_{ik_1 \dots k_{p-1}} Q_{i l_1 \dots l_{q-1}}$. This may sound abstract, but it's very much something you've seen before: given a pair of 1-tensors \mathbf{a} and \mathbf{b} , also known as vectors, we can combine them to get a 0-tensor, also known as a number

$$\mathbf{a} \cdot \mathbf{b} = a_i b_i$$

This, of course, is just the inner-product. It is a useful operation precisely because the 0-tensor on the right-hand side is, like all 0-tensors, independent of the choice of basis that we choose to express the vectors.

The Quotient Rule

In practice, it's not hard to recognise a tensor when you see one. In any setting, they're usually just objects with a bunch of i and j indices, each of which clearly transforms as a vector. If in doubt, you can just check explicitly how the thing transforms. (There are cases where this check is needed. In later [courses](#), you'll meet an object called the Levi-Civita connection Γ_{jk}^i which looks for all the world like a tensor but turns out, on closer inspection, to be something more subtle.)

There is a more formal way to say this. Let $T_{i_1 \dots i_{p+q}}$ be a bunch of numbers that you think might comprise a tensor of rank $p+q$ in some coordinate basis. If $T_{i_1 \dots i_{p+q}}$ are indeed the components of a tensor then you can feed it a rank q tensor $u_{j_1 \dots j_q}$ and it will spit back a rank p tensor

$$v_{i_1 \dots i_p} = T_{i_1 \dots i_p j_1 \dots j_q} u_{j_1 \dots j_q} \quad (6.5)$$

There is a converse to this statement. If for every tensor $u_{j_1 \dots j_q}$, the output $v_{i_1 \dots i_p}$ defined in (6.5) is a tensor, then $T_{i_1 \dots i_p j_1 \dots j_q}$ are the components of a tensor. This is called the *quotient rule*.

It is straightforward, if a little fiddly, to prove the quotient rule. It's sufficient to restrict attention to tensors u formed from the tensor product of vectors $u_{j_1 \dots j_q} = c_{j_1} \dots d_{j_q}$. Then, by assumption, $v_{i_1 \dots i_p} = T_{i_1 \dots i_p j_1 \dots j_q} u_{j_1 \dots j_q}$ is a tensor. If we then contract with p further vectors $\mathbf{a}, \dots, \mathbf{b}$ then $v_{i_1 \dots i_p} a_{i_1} \dots b_{i_p} = T_{i_1 \dots i_p j_1 \dots j_q} a_{i_1} \dots b_{i_p} c_{j_1} \dots d_{j_q}$ is necessarily a scalar. This is then enough to ensure the correct transformation rule (6.3) for the components $T_{i_1 \dots i_p j_1 \dots j_q}$.

Symmetry and Anti-Symmetry

The symmetrisation properties of tensors are worthy of comment. A tensor that obeys

$$T_{ijp\dots q} = \pm T_{jip\dots q}$$

is said to be *symmetric* (for $+$) or *anti-symmetric* (for $-$) in the indices i and j . If a tensor is (anti)-symmetric in one coordinate system then it is (anti)-symmetric in any coordinate system

$$T'_{ijp\dots q} = R_{ik} R_{jl} R_{pr} \dots R_{qs} T_{klr\dots s} = \pm R_{ik} R_{jl} R_{pr} \dots R_{qs} T_{lkr\dots s} = \pm T'_{jip\dots q}$$

A tensor that is (anti)-symmetric in all pairs of indices is said to be *totally (anti)-symmetric*. Note that for tensors in \mathbb{R}^n , there are no anti-symmetric tensors of rank $p > n$ because at least one of the indices must take the same value and so the tensor necessarily vanishes. A totally anti-symmetric tensor of rank p in \mathbb{R}^n has $\binom{n}{p}$ independent components.

Let's now restrict our attention to \mathbb{R}^3 . A tensor of rank 2 is our new fancy name for a 3×3 matrix T_{ij} . In general, it has 9 independent components. We can always decompose it into the symmetric and anti-symmetric pieces

$$S_{ij} = \frac{1}{2}(T_{ij} + T_{ji}) \quad \text{and} \quad A_{ij} = \frac{1}{2}(T_{ij} - T_{ji})$$

which have 6 and 3 independent components respectively. Our discussion above shows that S and A are each, themselves, tensors. In fact, the symmetric piece can be decomposed further,

$$S_{ij} = P_{ij} + \frac{Q}{3}\delta_{ij}$$

where $Q = S_{ii}$ is the trace of S and carries a single degree of freedom, while P_{ij} is the traceless part of S and carries 5. The importance of this decomposition is that A , P and Q are individually tensors. In contrast, if you were to take, say, the upper-left-hand component of the original matrix T_{ij} then that doesn't form a tensor.

In \mathbb{R}^3 , we can also rewrite an anti-symmetric matrix in terms of a vector,

$$A_{ij} = \epsilon_{ijk}B_k \quad \Longleftrightarrow \quad B_k = \frac{1}{2}\epsilon_{ijk}A_{ij}$$

The upshot is that in any 3×3 matrix can be decomposed as

$$T_{ij} = P_{ij} + \epsilon_{ijk}B_k + \frac{1}{3}\delta_{ij}Q \tag{6.6}$$

where $P_{ii} = 0$.

6.1.3 Invariant Tensors

There are two important invariant tensors in \mathbb{R}^n .

- We've met the first already: it is the rank 2 tensor δ_{ij} . As we noted previously, this is invariant because

$$\delta'_{ij} = R_{ik}R_{jl}\delta_{kl} = \delta_{ij}$$

Note that δ_{ij} is invariant under any $R \in O(n)$.

- The rank n totally anti-symmetric tensor $\epsilon_{i_1 \dots i_n}$. This is defined by $\epsilon_{12 \dots n} = +1$. If you swap any two indices you get a minus sign. In particular, if any two indices are repeated, the epsilon symbol vanishes. This is invariant because

$$\epsilon'_{i_1 \dots i_n} = R_{i_1 j_1} \dots R_{i_n j_n} \epsilon_{j_1 \dots j_n} = \det R \epsilon_{i_1 \dots i_n} = \epsilon_{i_1 \dots i_n}$$

Note that the epsilon symbol is only invariant under $R \in SO(n)$ but it is not invariant under $R \in O(n)$ with $\det R = -1$. It picks up a minus sign under reflections. The invariance of ϵ_{ijk} in \mathbb{R}^3 is the reason why the cross-product $(\mathbf{a} \times \mathbf{b})_i = \epsilon_{ijk} a_j b_k$ is itself a vector. Or, said differently, why the triple product $\mathbf{a} \cdot (\mathbf{b} \times \mathbf{c}) = \epsilon_{ijk} a_i b_j c_k$ is independent of the choice of basis.

In general, a tensor is said to be invariant under a given rotation R if

$$T'_{i_1 \dots i_n} = R_{i_1 j_1} \dots R_{i_n j_n} T_{j_1 \dots j_n} = T_{i_1 \dots i_n}$$

A tensor that is invariant under all rotations R is said to be *isotropic*. Obviously all tensors of rank 0 are isotropic. What about higher rank tensors?

Claim: The only non-zero isotropic tensors in \mathbb{R}^3 of rank $p = 1, 2$ or 3 are $T_{ij} = \alpha \delta_{ij}$ and $T_{ijk} = \beta \epsilon_{ijk}$ with α and β constant. In particular, there are no isotropic tensors of rank 1 (essentially because a vector always points in a preferred direction).

Proof: The idea is simply to look at how tensors transform under a bunch of specific rotations by π or $\pi/2$ about certain axes.

For example, consider a tensor of rank 1, so that

$$T'_i = R_{ij} T_j \quad \text{with} \quad R_{ij} = \begin{pmatrix} -1 & 0 & 0 \\ 0 & -1 & 0 \\ 0 & 0 & +1 \end{pmatrix} \quad (6.7)$$

Requiring $T'_i = T_i$ gives $T_1 = T_2 = 0$. Clearly a similar argument, using a different R , also gives $T_3 = 0$.

For a tensor of rank 2, consider the transformation

$$T'_{ij} = \tilde{R}_{ik} \tilde{R}_{jl} T_{kl} \quad \text{with} \quad \tilde{R}_{ij} = \begin{pmatrix} 0 & 1 & 0 \\ -1 & 0 & 0 \\ 0 & 0 & +1 \end{pmatrix} \quad (6.8)$$

which is a rotation by $\pi/2$ about the z -axis. The rotation gives $T'_{13} = T_{23}$ and $T'_{23} = -T_{13}$ so if $T'_{ij} = T_{ij}$, we must have $T_{13} = T_{23} = 0$. Meanwhile $T'_{11} = T_{22}$. Similar arguments tell us that all off-diagonal elements must vanish and all diagonal elements must be equal: $T_{11} = T_{22} = T_{33} = \alpha$ for some α . Hence $T_{ij} = \alpha \delta_{ij}$.

Finally, for a rank 3 tensor we have

$$T'_{ijk} = R_{il}R_{jp}R_{kq}T_{lpq}$$

If we pick R given in (6.7), then we find $T'_{133} = -T_{133}$ and $T'_{111} = -T_{111}$. Similar arguments show that an isotropic tensor must have $T_{ijk} = 0$ unless i, j and k are all distinct. Meanwhile, if we pick $R = \tilde{R}$ given in (6.8), then we get $T'_{123} = -T_{213}$. We end up with the result we wanted: T_{ijk} is isotropic if and only if $T_{ijk} = \beta\epsilon_{ijk}$ for some constant β . \square

Although we won't prove it here, all other isotropic tensors can be formed from δ_{ij} and ϵ_{ijk} . For example, the only isotropic 4-tensor in \mathbb{R}^3 is

$$T_{ijkl} = \alpha\delta_{ij}\delta_{kl} + \beta\delta_{ik}\delta_{jl} + \gamma\delta_{il}\delta_{jk}$$

with α, β and γ constants. You could try to cook up something involving ϵ_{ijk} but it doesn't give anything new. In particular, $\epsilon_{ijk}\epsilon_{ilp} = \delta_{jl}\delta_{kp} - \delta_{jp}\delta_{kl}$.

There is also an analogous result in \mathbb{R}^n : all isotropic tensors can be constructed from the symmetric 2-tensor δ_{ij} and the totally anti-symmetric n -tensor $\epsilon_{i_1\dots i_n}$.

Invariant Integrals

It is sometimes possible to use invariance properties to immediately write down the index structure of an integral, without doing the hard work of evaluating everything term by term. Suppose that we have some integral of the form

$$T_{ij\dots k} = \int_V f(r)x_ix_j\dots x_k dV$$

with $r = |\mathbf{x}|$. Then under a rotation, we have

$$T'_{ij\dots k} = R_{ip}R_{jq}\dots R_{kr}T_{pq\dots r} = \int_V f(r)x'_ix'_j\dots x'_k dV$$

with, as usual, $x'_i = R_{ij}x_j$. But if we now change the integration variables to x' , both $r = |\mathbf{x}| = |\mathbf{x}'|$ and $dV = dV'$ are invariant. (The latter because the Jacobian is $\det R = 1$). If the domain of integration is also rotationally invariant, so $V = V'$, then the final result must itself be an invariant tensor, $T'_{ij\dots k} = T_{ij\dots k}$.

Here are some examples. First, suppose that we have a 3d integral over the interior of a sphere of radius R , given by

$$T_i = \int_V \rho(r) x_i dV \quad (6.9)$$

This must be equal to some invariant 1-tensor (i.e. a vector), but there are no such objects. In other words, we can say immediately that $T_i = 0$. You can check this straightforwardly by doing the integral in, say, spherical polar coordinates.

Things change if we look at an integral with two hanging indices,

$$T_{ij} = \int_V \rho(r) x_i x_j dV \quad (6.10)$$

(In Section 6.2, we will find integrals of this form arising when we compute the inertia tensor of a sphere.) By the argument above T_{ij} must be an isotropic tensor and hence proportional to δ_{ij} ,

$$T_{ij} = \int_V \rho(r) x_i x_j dV = \alpha \delta_{ij}$$

for some α . If we take the trace, we get

$$\int_V \rho(r) r^2 dV = 3\alpha$$

Hence,

$$T_{ij} = \frac{1}{3} \delta_{ij} \int_V \rho(r) r^2 dV = \frac{4\pi}{3} \delta_{ij} \int_0^R dr \rho(r) r^4 \quad (6.11)$$

For example, if $\rho(r) = \rho_0$ is constant, then $T_{ij} = \frac{4}{15} \pi \rho_0 R^5 \delta_{ij}$.

Here's a slightly more complicated example (taken from the calculation of Stokes flow around a sphere in [Fluid Mechanics](#)). Consider the surface integral over a sphere of radius R ,

$$\tilde{T}_k = a_j \int_{\mathbf{S}^2} dS_i \frac{x_i x_j x_k}{r^5}$$

This time we have a vector \mathbf{a} in the game, so it must be the case that $\tilde{T}_k = \beta a_k$ for some constant β . One way to compute β is to strip off the vector \mathbf{a} and instead look at

$$\tilde{T}_{jk} = \int_{\mathbf{S}^2} dS_i \frac{x_i x_j x_k}{r^5} = \beta \delta_{jk}$$

which now should be proportional to the invariant tensor δ_{jk} as shown, with the same coefficient β since $\tilde{T}_k = T_{jk}a_j = \beta a_k$. At this point, we again take the trace over the j and k indices to get

$$\int_{\mathbf{S}^2} dS_i \frac{x_i x_j x_j}{r^5} = 3\beta$$

But this integral is given by

$$\int_{\mathbf{S}^2} dS_i \frac{x_i}{r^3} = \int_{\mathbf{S}^2} d\mathbf{S} \cdot \frac{\mathbf{n}}{r^2} = 4\pi$$

and so we get $\beta = 4\pi/3$.

6.1.4 Tensor Fields

A tensor field over \mathbb{R}^3 is the assignment of a tensor $T_{i\dots k}(\mathbf{x})$ to every point $\mathbf{x} \in \mathbb{R}^3$. This is the generalisation of a vector field

$$\mathbf{F} : \mathbb{R}^3 \rightarrow \mathbb{R}^3$$

to a map of the kind

$$T : \mathbb{R}^3 \rightarrow \mathbb{R}^m$$

with m the number of components of the tensor. So, for example, a map that assigns a symmetric, traceless rank 2 tensor $P_{ij}(\mathbf{x})$ to every point has $m = 5$.

The tensor field $T_{i\dots k}(\mathbf{x})$ is sometimes denoted as $T_{i\dots k}(x^l)$ which is supposed to show that the field depends on all coordinates x^1, \dots, x^3 . It's not great notation because the indices as subscripts are supposed to take some definite values, while the index l in the argument is supposed to denote the whole set of indices. It's especially bad notation when combined with the summation convention and we won't adopt it here.

There is one very famous example of a tensor field. Einstein's theory of general relativity is described by a rank 2 tensor at every point in space. This is called the *metric*. The dynamics of this rank 2 tensor field describe gravity. (I've brushed something rather important under the rug here. Einstein's theory is a rank 2 tensor in *spacetime*, not just in space. Which means that the rank 2 tensor is a 4×4 matrix, rather than a 3×3 matrix.)

Before we move on, it's worth pausing to mention a slightly subtle point. Not all maps $\mathbb{R}^3 \rightarrow \mathbb{R}^3$ qualify as "vector fields". The point \mathbf{x} in the codomain \mathbb{R}^3 is a vector and so its components transform in the appropriate way under rotation. To be a vector field, the components of the map must transform under the *same* rotation. Similar comments hold for a tensor field.

To illustrate this, the electric field $\mathbf{E}(\mathbf{x})$ is an example of a vector field. If you rotate in space, and so change \mathbf{x} , then the direction \mathbf{E} also changes: the rotation acts on both the argument \mathbf{x} and the function itself \mathbf{E} .

In contrast, there are maps $\mathbb{R}^3 \rightarrow \mathbb{R}^3$ where, although the domain and codomain have the same dimension, vectors in them transform under different rotations. For example, in particle physics there exists an object called a *quark field* which, for our (admittedly, slightly dumbed down) purposes, can be thought of as a map $\mathbb{R}^3 \rightarrow \mathbb{R}^3$. This is a quantum field whose ripples are the particles that we call quarks, but these details can be safely ignored for the next couple of years of your life. We will write this field as $q_a(\mathbf{x})$ where the $a = 1, 2, 3$ label is the “colour” of the quark. If we rotate in space, then \mathbf{x} changes but the colour of the quark does not. There is then an independent rotation that acts on the codomain and rotates the colour, but leaves the point in space unchanged. Because the rotations that act on the domain and codomain are unrelated, the quark field is usually not referred to as a vector field.

Taking Derivatives

Given a tensor field, we can always construct higher rank tensors by taking derivatives. In fact, we’ve already seen a prominent example of this earlier in these lectures. There, we started with a scalar field $\phi(\mathbf{x})$ and differentiated to get the gradient $\nabla\phi$. This means that we start with a rank 0 tensor and differentiate to get a rank 1 tensor.

Strictly speaking, we didn’t previously prove that $\nabla\phi$ is a vector field. But it’s straightforward to do so. As we’ve seen above, we need to show that it transforms correctly under rotations. Any vector \mathbf{v} can be decomposed in two different ways,

$$\mathbf{v} = v^i \mathbf{e}_i = v'^i \mathbf{e}'_i$$

where $\{\mathbf{e}_i\}$ and $\{\mathbf{e}'_i\}$ are two orthonormal bases, each obeying $\mathbf{e}_i \cdot \mathbf{e}_j = \mathbf{e}'_i \cdot \mathbf{e}'_j = \delta_{ij}$, and v^i and v'^i are the two different coordinates for \mathbf{v} . If we expand \mathbf{x} in this way

$$\mathbf{x} = x_i \mathbf{e}_i = x'_i \mathbf{e}'_i \implies x_i = (\mathbf{e}_i \cdot \mathbf{e}'_j) x'_j \implies \frac{\partial x^i}{\partial x'^j} = \mathbf{e}_i \cdot \mathbf{e}'_j$$

Here $\mathbf{e}_i \cdot \mathbf{e}'_j$ is the rotation matrix that takes us from one basis to the other. Meanwhile, we can always expand one set of basis vectors in terms of the other,

$$\mathbf{e}_i = (\mathbf{e}_i \cdot \mathbf{e}'_j) \mathbf{e}'_j = \frac{\partial x^i}{\partial x'^j} \mathbf{e}'_j$$

This tells us that we could equally as well write the gradient as

$$\nabla\phi = \frac{\partial\phi}{\partial x^i} \mathbf{e}_i = \frac{\partial\phi}{\partial x^i} \frac{\partial x^i}{\partial x'^j} \mathbf{e}'_j = \frac{\partial\phi}{\partial x'^j} \mathbf{e}'_j$$

This is the expected result: if you work in a different primed basis, then you have the same definition of $\nabla\phi$, but just with primes on both \mathbf{e}'_i and $\partial/\partial x'^i$. This means that the components $\partial_i\phi$ transform correctly under a rotation, so $\nabla\phi$ is indeed a vector.

We can extend the result above to any, suitably smooth, tensor field $T(\mathbf{x})$ of rank p . We can differentiate this any number of times to get a new tensor field of rank, say, $p + q$,

$$X_{i_1 \dots i_q j_1 \dots j_p} = \frac{\partial}{\partial x_{i_1}} \cdots \frac{\partial}{\partial x_{i_q}} T_{j_1 \dots j_p}(\mathbf{x}) \quad (6.12)$$

To verify that this is indeed a tensor, we need to check how it changes under a rotation. In a new basis, we have $x'_i = R_{ij}x_j$ (where $R_{ij} = \mathbf{e}'_i \cdot \mathbf{e}_j$ in the notation above) and so

$$\frac{\partial x'_i}{\partial x_j} = R_{ij} \implies \frac{\partial}{\partial x'_i} = \frac{\partial x_j}{\partial x'_i} \frac{\partial}{\partial x_j} = R_{ij} \frac{\partial}{\partial x_j}$$

which is the result we need for X in (6.12) to qualify as a tensor field.

We can implement any of the tensorial manipulations that we met previously for tensor fields. For example, if we start with a vector field $\mathbf{F}(\mathbf{x})$, we can form a rank 2 tensor field

$$T_{ij}(\mathbf{x}) = \frac{\partial F_i}{\partial x_j}$$

But we saw in (6.6) that any rank 2 tensor field can be decomposed into various pieces. There is an anti-symmetric piece

$$A_{ij}(\mathbf{x}) = \epsilon_{ijk} B_k(\mathbf{x}) \quad \text{with} \quad B_k = \frac{1}{2} \epsilon_{ijk} \frac{\partial F_i}{\partial x_j} = -\frac{1}{2} (\nabla \times \mathbf{F})_k$$

and a trace piece

$$Q = \frac{\partial F_i}{\partial x_i} = \nabla \cdot \mathbf{F}$$

and, finally, a symmetric, traceless piece

$$P_{ij}(\mathbf{x}) = \frac{1}{2} \left(\frac{\partial F_i}{\partial x_j} + \frac{\partial F_j}{\partial x_i} \right) - \frac{1}{3} \nabla \cdot \mathbf{F}$$

Obviously, the first two of these are familiar tensors (in this case a scalar and vector) from earlier sections.

6.2 Physical Examples

Our discussion above was rooted firmly in mathematics. There are many places in physics where tensors appear. Here we give a handful of examples.

6.2.1 Electric Fields in Matter

Apply an electric field \mathbf{E} to a lump of stuff. A number of things can happen.

If the lump of stuff is an insulator then the material will become *polarised*. This means that the positive electric charge will be pushed in one direction, the negative in another until the lump of stuff acts like a dipole. (This is described in some detail in Section 7 of the lectures on [Electromagnetism](#).) One might think that the resulting polarisation vector \mathbf{P} points in the same direction as the electric field \mathbf{E} , but that's too simplistic. For many lumps of stuff, the underlying crystal structure allows the electric charges to shift more freely in some directions than others. The upshot is that the relation between polarisation \mathbf{P} and applied electric field \mathbf{E} is given by

$$\mathbf{P} = \alpha \mathbf{E}$$

where α is a matrix known as the *polarisation tensor*. In a given basis, it has components α_{ij} .

There is a similar story if the lump of stuff is a conductor. This time an applied electric field gives rise to a current density \mathbf{J} . Again, the current is not necessarily parallel to the electric field. The relationship between them is now

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

This is known as *Ohm's law*. In general σ is a 3×3 matrix known as the *conductivity tensor*; in a given basis, it has components σ_{ij} .

What can we say about σ when the material is isotropic, meaning that it looks the same in all directions? In this case, no direction is any different from any other. With no preferred direction, the conductivity tensor must be proportional to an invariant tensor, so that it looks the same in all coordinate systems. What are our options?

For 3d materials, the only option is $\sigma_{ij} = \sigma \delta_{ij}$, which ensures that the current does indeed run parallel to the electric field. In this case σ is just referred to as the *conductivity*.

However, suppose that we're dealing with a thin wafer of material in which both the current and electric field are restricted to lie in a plane. This changes the story because now we're dealing with vectors in \mathbb{R}^2 rather than \mathbb{R}^3 and \mathbb{R}^2 is special because there are two invariant 2-tensors in this dimension: δ_{ij} and ϵ_{ij} . This means that the most general conductivity tensor for an isotropic 2d material takes the form

$$\sigma_{ij} = \sigma_{xx}\delta_{ij} + \sigma_{xy}\epsilon_{ij} = \begin{pmatrix} \sigma_{xx} & \sigma_{xy} \\ -\sigma_{xy} & \sigma_{xx} \end{pmatrix}$$

Here σ_{xx} is called the *longitudinal conductivity* while σ_{xy} is called the *Hall conductivity*. If $\sigma_{xy} \neq 0$ then an electric field in the x -direction induces a current in the y -direction.

As an aside, it turns out that the seemingly mundane question of understanding σ_{xy} in real materials is closely tied to some of the most interesting breakthroughs in mathematics in recent decades! This is the subject of the [Quantum Hall Effect](#).

6.2.2 The Inertia Tensor

Another simple example of a tensor arises in Newtonian mechanics. A *rigid body* rotating about the origin can be modelled by some number of masses m_a at positions \mathbf{x}_a , all moving with velocity $\dot{\mathbf{x}}_a = \boldsymbol{\omega} \times \mathbf{x}_a$. Here $\boldsymbol{\omega}$ is known as the angular velocity. The angular velocity $\boldsymbol{\omega}$ is related to the angular momentum \mathbf{L} by

$$\mathbf{L} = I\boldsymbol{\omega} \tag{6.13}$$

with I the *inertia tensor*. The angular momentum does not necessarily lie parallel to the angular velocity and, correspondingly, I is in general a matrix, rather than a single number. In fact, we can easily derive an expression for the inertia tensor. The angular momentum is

$$\mathbf{L} = \sum_a m_a \mathbf{x}_a \times \dot{\mathbf{x}}_a = \sum_a m_a \mathbf{x}_a \times (\boldsymbol{\omega} \times \mathbf{x}_a) = \sum_a m_a \left(|\mathbf{x}_a|^2 \boldsymbol{\omega} - (\mathbf{x}_a \cdot \boldsymbol{\omega}) \mathbf{x}_a \right)$$

In components, $L_i = I_{ij}\omega_j$, where

$$I_{ij} = \sum_a m_a \left(|\mathbf{x}_a|^2 \delta_{ij} - (\mathbf{x}_a)_i (\mathbf{x}_a)_j \right)$$

For a continuous object with density $\rho(\mathbf{x})$, we can replace the sum with a volume integral

$$I_{ij} = \int_V \rho(\mathbf{x}) \left(|\mathbf{x}|^2 \delta_{ij} - x_i x_j \right) dV \tag{6.14}$$

So, for example, $I_{33} = \int \rho(x_1^2 + x_2^2) dV$ and $I_{12} = -\int \rho x_1 x_2 dV$.

An Example: A Sphere

For a ball of radius R and density $\rho(r)$, the inertia tensor is

$$I_{ij} = \int_V \rho(r)(r^2\delta_{ij} - x_i x_j) dV$$

The second of these terms is the integral (6.10) that we simplified in Section 6.1.3 using isotropy arguments. Using (6.11), we have

$$I_{ij} = \frac{2}{3}\delta_{ij} \int_V \rho(r)r^2 dV = \frac{8\pi}{3}\delta_{ij} \int_0^R dr \rho(r)r^4$$

For example, if $\rho(r) = \rho_0$ is constant, then $I_{ij} = \frac{8}{15}\pi\rho_0 R^5\delta_{ij} = \frac{2}{5}MR^2\delta_{ij}$ where M is the mass of the sphere.

Another Example: A Cylinder

The sphere is rather special because the inertia tensor is proportional to δ_{ij} . That's not the case more generally. Consider, for example, a solid 3d cylinder of radius a and height $2L$, with uniform density ρ . The mass is $M = 2\pi a^2 L \rho$. We align the cylinder with the z -axis and work in cylindrical polar coordinates $x = r \cos \phi$ and $y = r \sin \phi$. The components of the inertia tensor are then

$$\begin{aligned} I_{33} &= \int_V \rho(x^2 + y^2) dV = \rho \int_0^{2\pi} d\phi \int_0^a dr \int_{-L}^{+L} dz r r^2 = \rho\pi L a^4 \\ I_{11} &= \int_V \rho(y^2 + z^2) dV = \rho \int_0^{2\pi} d\phi \int_0^a dr \int_{-L}^{+L} dz r(r^2 \sin^2 \phi + z^2) = \rho\pi a^2 L \left(\frac{a^2}{2} + \frac{2L^2}{3} \right) \end{aligned}$$

By symmetry, $I_{22} = I_{11}$. For the off-diagonal elements, we have

$$I_{13} = - \int_V \rho x_1 x_3 dV = -\rho \int_0^{2\pi} d\phi \int_0^a dr \int_{-L}^L dz r^2 z \cos \phi = 0$$

where the integral vanishes due to the ϕ integration. Similarly, $I_{12} = I_{13} = 0$. We find that the inertia tensor for the cylinder is

$$I = \text{diag} \left(M \left(\frac{a^2}{4} + \frac{L^2}{3} \right), M \left(\frac{a^2}{4} + \frac{L^2}{3} \right), \frac{1}{2} M a^2 \right) \quad (6.15)$$

Note that the inertia tensor is diagonal in our chosen coordinates.

The Eigenvectors of the Inertia Tensor

The inertia tensor I defined in (6.14) has a special property: it is symmetric

$$I_{ij} = I_{ji}$$

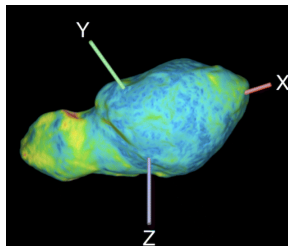
Any symmetric matrix I can always be diagonalised by an appropriate rotation. This means that there exists an $R \in SO(n)$ such that

$$I' = RIR^T = \text{diag}(I_1, I_2, I_3)$$

Another way of saying this is that any symmetric rank 2 tensor has a basis of orthonormal eigenvectors $\{\mathbf{e}_i\}$, with I_i the corresponding eigenvalues.

In the case of the inertia tensor, the eigenvectors \mathbf{e}_1 , \mathbf{e}_2 and \mathbf{e}_3 are called the *principal axes* of the solid. It means that any object, no matter how complicated, has its own preferred set of orthonormal axes embedded within it. If the object has some symmetry, then the principal axes will always be aligned with this symmetry. This, for example, was the case for the cylinder that we computed above where aligning the cylinder with the z -axis automatically gave us a diagonal inertia tensor (6.15).

In general, it will be less obvious where the principal axes lie. For example, the figure on the right shows the asteroid Toutatis, which is notable for its lumpy shape. The principal axes are shown embedded in the asteroid.



From (6.13), the angular momentum \mathbf{L} is aligned with the angular velocity $\boldsymbol{\omega}$ only if a body spins about one of its principal axes. It turns out that, in this case, nice things happen and the body spins smoothly.

However, if \mathbf{L} and $\boldsymbol{\omega}$ are misaligned, the body exhibits more complicated tumbling, wobbling motion as it spins. You can learn all about this in the lectures on [Classical Dynamics](#). (For what it's worth, Toutatis does not spin about a principal axes.)

6.2.3 Higher Rank Tensors

You might reasonably complain that, after all that work defining tensors, the examples that we've given here are nothing more exotic than matrices, mapping one vector to another. And you would be right. However, as we get to more sophisticated theories of physics, tensors of higher rank do make an appearance. Here we don't give full details, but just say a few words to give you a flavour of things to come.

Perhaps the simplest example arises in the theory of elastic materials. These materials can be subjected to *strain*, which describes the displacement of the material at each point, and *stress*, which describes the forces acting on the material at each point. But each of these is itself a 2-tensor (strictly a tensor field). The *strain tensor* e_{ij} is a symmetric tensor that describes the way the displacement in the x^i direction varies in the x^j . The *stress tensor* σ_{ij} describes the component of the force F_i across a plane normal to x^j . These two tensors are related by

$$\sigma_{ij} = C_{ijkl}e_{kl}$$

This is the grown up version of Hooke's law. In general an elastic material is characterised by the *elasticity tensor*, also known as the *stiffness tensor*, C_{ijkl} .

Higher rank tensors also appear prominently in more advanced descriptions of geometry. In higher dimensions, the simple Gaussian curvature that we met in Section 2 isn't enough to capture all the interesting ways in which spaces can curve in different directions. Instead, it is replaced by a 4-tensor R_{ijkl} known as the Riemann curvature. In the context of physics, this 4-tensor describes the bending of space and time and is needed for the grown-up version of Newton's law of gravity.

6.3 A Unification of Integration Theorems

In this final section, we turn back to matters of mathematics. The three integral theorems that we met in Section 4 are obviously closely related. To end these lectures, we show how they can be presented in a unified framework. This requires us to introduce some novel and slightly formal ideas. These go quite a bit beyond what is usually covered in an introductory course on vector calculus, but we will meet these objects again in later courses on [Differential Geometry and General Relativity](#). View this section as a taste of things to come.

6.3.1 Integrating in Higher Dimensions

Our unified framework will give us integral theorems in any dimension \mathbb{R}^n . If you look back at Section 4, you'll notice that the divergence theorem already holds in any \mathbb{R}^n . Meanwhile, Stokes' theorem is restricted to surfaces in \mathbb{R}^3 for the very simple reason that the cross-product is only defined in \mathbb{R}^3 . This suggests that before we can extend our integral theorems to higher dimensions, we should first ask a more basic question: how do we extend the cross product to higher dimensions?

The introduction of tensors gives us a way to do this. Given two vectors \mathbf{a} and \mathbf{b} in \mathbb{R}^3 , the cross-product is

$$(\mathbf{a} \times \mathbf{b})_i = \epsilon_{ijk} a_j b_k$$

From this perspective, the reason that the cross product can only be employed in \mathbb{R}^3 is because it's only there that the ϵ_{ijk} symbol has three entries. If, in contrast, we're in \mathbb{R}^4 then we have ϵ_{ijkl} and so if we feed it two vectors \mathbf{a} and \mathbf{b} , then we find ourselves with a tensor of rank 2, $T_{ij} = \epsilon_{ijkl} a_k b_l$.

The tensors that we get from an epsilon symbol are always special, in the sense that they are totally anti-symmetric. The anti-symmetry condition doesn't impose any extra constraint on a 0-tensor ϕ or a 1-tensor a_i as these are just scalar fields and vector fields respectively. It only kicks in when we get to tensors of rank 2 or higher.

With this in mind, we can revisit the cross product. We can define the cross product in any dimension \mathbb{R}^n : it is a map that eats two vectors \mathbf{a} and \mathbf{b} and spits back an anti-symmetric $(n - 2)$ -tensor

$$(\mathbf{a} \times \mathbf{b})_{i_1 \dots i_{n-2}} = \epsilon_{i_1 \dots i_n} a_{i_{n-1}} b_{i_n}$$

The only thing that's special about \mathbb{R}^3 is that we get back another vector, rather than a higher dimensional tensor.

There is also a slightly different role played by the epsilon symbol $\epsilon_{i_1, \dots, i_n}$: it provides a map from anti-symmetric p -tensors to anti-symmetric $(n - p)$ -tensors, simply by contracting indices,

$$\epsilon : T_{i_1 \dots i_p} \mapsto \frac{1}{(n - p)!} \epsilon_{i_1 \dots i_n} T_{i_{n-p+1} \dots i_n} \quad (6.16)$$

This map goes by the fancy name of the *Hodge dual*. (Actually, it's an entirely trivial version of the Hodge dual. The proper Hodge dual is a generalisation of this idea to curved spaces.)

Our next step is to think about what this has to do with integration. Recall that earlier in these lectures we found two natural ways to integrate vector fields in \mathbb{R}^3 . The first is along a line

$$\int_C \mathbf{F} \cdot d\mathbf{x} \quad (6.17)$$

which captures the component vector field *tangent* to the line. We can perform this procedure in any dimension \mathbb{R}^n . The second operation is to integrate a vector field over a surface

$$\int_S \mathbf{F} \cdot d\mathbf{S} \quad (6.18)$$

where $d\mathbf{S}$ points in the direction normal to the surface. This integration captures the component of the vector field *normal* to the surface and only makes sense in \mathbb{R}^3 . This is because it's only in \mathbb{R}^3 that a two-dimensional surface has a unique normal. More operationally, this normal, which is buried in the definition of $d\mathbf{S}$, requires us to use the cross product. For a parameterised surface $\mathbf{x}(u, v)$, the vector area element is

$$d\mathbf{S} = \frac{\partial \mathbf{x}}{\partial u} \times \frac{\partial \mathbf{x}}{\partial v} du dv$$

or, in components,

$$dS_i = \epsilon_{ijk} \frac{\partial x^j}{\partial u} \frac{\partial x^k}{\partial v} du dv$$

Now comes a mathematical sleight of hand. Rather than thinking of (6.18) as the integral of a vector field projected normal to the surface, instead think of it as the integral of an anti-symmetric 2-tensor $F_{ij} = \epsilon_{ijk} F_k$ integrated *tangent* to the surface. We then have

$$\int_S \mathbf{F} \cdot d\mathbf{S} = \int_S F_{ij} dS_{ij} \quad \text{with} \quad dS_{ij} = \frac{1}{2} \left(\frac{\partial x^j}{\partial u} \frac{\partial x^k}{\partial v} - \frac{\partial x^j}{\partial v} \frac{\partial x^k}{\partial u} \right) du dv \quad (6.19)$$

This is the same equation as before, just with the epsilon symbol viewed as part of the integrand F_{ij} rather than as part of the measure dS_i . Note that we've retained the anti-symmetry of the area element dS_{ij} that was inherent in our original cross product definition of $d\mathbf{S}$. Strictly speaking this isn't necessary because we're contracting with anti-symmetric indices in F_{ij} , but it turns out that it's best to think of both objects F_{ij} and dS_{ij} as individually anti-symmetric.

This new perspective suggests a way to generalise to higher dimensions. In the line integral (6.17) we're integrating a vector field over a line. In the surface integral (6.19), we're really integrating an anti-symmetric 2-tensor over a surface. The key idea is that one can integrate a totally anti-symmetric p -tensor over a p -dimensional subspace.

Specifically, given an anti-symmetric p -tensor, the generalisation of the line integral (6.17) is the integration over a p -dimensional subspace,

$$\int_M T_{i_1 \dots i_p} dS_{i_1 \dots i_p} \quad (6.20)$$

where $\dim(M) = p$. Here $dS_{i_1 \dots i_p}$ is a higher dimensional version of the “area element” defined in (6.19).

Alternatively, the higher dimensional version of the surface integral (6.18) involves first mapping the p -tensor to an $(n - p)$ -tensor using the Hodge dual. This can subsequently be integrated over an $(n - p)$ -dimensional subspace

$$\int_{\tilde{M}} T_{i_1 \dots i_p} \epsilon_{i_1 \dots i_p j_1 \dots j_{n-p}} d\tilde{S}_{j_1 \dots j_{n-p}} \quad (6.21)$$

with $\dim(\tilde{M}) = n - p$.

In fact, we’ve already met an integral of the form (6.21) elsewhere in these lectures, since this is what we’re implicitly doing when we integrate a scalar field over a volume. In this case the “area element” is just $dS_{i_1 \dots i_n} = \frac{1}{n!} \epsilon_{i_1 \dots i_n} dV$ and the two epsilon symbols just multiply to a constant.. When actually computing a volume integral, this extra machinery is more of a distraction than a help.. But if we want to know how to think about things more generally then it’s extremely useful.

6.3.2 Differentiating Anti-Symmetric Tensors

We’ve now learned how to integrate anti-symmetric tensors. Our next step is to learn how to differentiate them. We’ve already noted in (6.12) that we can differentiate a p tensor once to get a tensor of rank $p + 1$, but in general differentiating loses the anti-symmetry property. As we now explain, there is a way to restore it so that when we differentiate a totally anti-symmetric p tensor, we end up with a totally anti-symmetric $(p + 1)$ -tensor.

For a scalar field, things are trivial. We can construct a vector field $\nabla\phi$ and this is automatically “anti-symmetric” because there’s nothing to anti-symmetrise.

If we’re given a vector field F_i , we can differentiate and then anti-symmetrise by hand. I will introduce a new symbol for “differentiation and anti-symmetrisation” and write

$$(\mathcal{D}F)_{ij} := \frac{1}{2} \left(\frac{\partial F_i}{\partial x^j} - \frac{\partial F_j}{\partial x^i} \right)$$

where the anti-symmetry is manifest on the right-hand side. I should confess that the notation $\mathcal{D}F$ is not at all standard. In subsequent courses, this object is usually viewed as something called a “differential form” and written simply as dF but the notation dF is loaded with all sorts of other connotations which are best ignored at this stage. Hence the made-up notation $\mathcal{D}F$.

In \mathbb{R}^3 , this anti-symmetric differentiation is equivalent to the curl using the Hodge map (6.16),

$$(\nabla \times \mathbf{F})_i = \epsilon_{ijk}(\mathcal{D}F)_{jk}$$

But now we can extend this definition to any anti-symmetric p -tensor. We can always differentiate and anti-symmetrise to get a $(p+1)$ -tensor defined by

$$(\mathcal{D}T)_{i_1 \dots i_{p+1}} = \frac{1}{p+1} \left(\frac{\partial T_{i_1 \dots i_p}}{\partial x^{i_{p+1}}} + p \text{ further terms} \right)$$

where the further terms involve replacing the derivative $\partial/\partial x^{i_{p+1}}$ with one of the other coordinates $\partial/\partial x^j$ so that the whole shebang is fully anti-symmetric.

Note that, with this definition of \mathcal{D} , if we differentiate twice then we take a p -tensor to a $(p+2)$ -tensor. But this $(p+2)$ -tensor always vanishes!

$$(\mathcal{D}\mathcal{D}T)_{i_1 \dots i_{p+2}} = 0$$

for any tensor T . This is because we’ll have two derivatives contracted with an epsilon and is the higher dimensional generalisation of the statements that $\nabla \times \nabla \phi = 0$ or $\nabla \cdot (\nabla \times \mathbf{F}) = 0$.

As an aside: this is actually the second time in these lectures that we’ve seen something vanish when you act twice, although you’d be forgiven for failing to notice the connection. Here our new anti-symmetric derivative obeys $\mathcal{D}^2(\text{anything}) = 0$. But we previously saw that the “boundary of a boundary” is always zero. This means that if a higher dimensional space (really a manifold) M has boundary ∂M then $\partial(\partial M) = 0$. Conceptually, these two ideas are very different but one can’t help but be struck by the similarity of the equations $\mathcal{D}^2(\text{anything}) = 0$ and $\partial^2(\text{anything}) = 0$, even though the “anything”’s are very different objects in the two formulae. It turns out that this similarity is pointing at a deep connection between the topology of spaces and the kinds of tensors that one can put on these spaces. In fancy maths words, this is the link between homology and cohomology.

Finally, we can now state the general integration theorem. Given an anti-symmetric p -tensor T , then

$$\int_M (\mathcal{D}T)_{i_1 \dots i_{p+1}} dS_{i_1 \dots i_{p+1}} = \int_{\partial M} T_{i_1 \dots i_p} dS_{i_1 \dots i_p} \quad (6.22)$$

Here $\dim(M) = p + 1$ and, therefore the boundary has $\dim(\partial M) = p$. Note that we don't use a different letter to distinguish the integration measure over these various spaces: everything is simply dS and you have to look closer at the indices to see what kind of space you're integrating over.

The equation (6.22) is a unification of all integration theorems. It contains the fundamental theorem of calculus (when $p = 0$), the divergence theorem (when $p = n - 1$) and Stokes' theorem (when $p = 1$ and $\mathbb{R}^n = \mathbb{R}^3$). Geometers refer to this generalised theorem simply as *Stokes' theorem* since that is the original result that it resembles most. The proof is simply a higher dimensional version of the proofs that we sketched previously.

There is, to put it mildly, quite a lot that I'm sweeping under the rug in the discussion above. In particular, the full Stokes' theorem does not hold only in \mathbb{R}^n but in a general curved space known as a manifold. In that context, one has to be a lot more careful about what kind of tensors we're dealing with and, as I mentioned above, Stokes' theorem should be written using a kind of anti-symmetric tensor known as a *differential form*. None of this really matters when working in flat space, but the differences become crucial when thinking about curved spaces. If you want to learn more, these topics will be covered in glorious detail in later courses on Differential Geometry or, for physicists, [General Relativity](#).

What You Really Need

Here are expressions for div, grad, curl and the Laplacian in various coordinate systems.

Cartesian: $\mathbf{x} = (x, y, z)$

$$\nabla f = \frac{\partial f}{\partial x} \hat{\mathbf{x}} + \frac{\partial f}{\partial y} \hat{\mathbf{y}} + \frac{\partial f}{\partial z} \hat{\mathbf{z}}$$

$$\nabla \cdot \mathbf{F} = \frac{\partial F_x}{\partial x} + \frac{\partial F_y}{\partial y} + \frac{\partial F_z}{\partial z}$$

$$\nabla \times \mathbf{F} = \left(\frac{\partial F_z}{\partial y} - \frac{\partial F_y}{\partial z} \right) \hat{\mathbf{x}} + \left(\frac{\partial F_x}{\partial z} - \frac{\partial F_z}{\partial x} \right) \hat{\mathbf{y}} + \left(\frac{\partial F_y}{\partial x} - \frac{\partial F_x}{\partial y} \right) \hat{\mathbf{z}}$$

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

Cylindrical Polars: $\mathbf{x} = (\rho \cos \phi, \rho \sin \phi, z)$

$$\nabla f = \frac{\partial f}{\partial \rho} \hat{\boldsymbol{\rho}} + \frac{1}{\rho} \frac{\partial f}{\partial \phi} \hat{\boldsymbol{\phi}} + \frac{\partial f}{\partial z} \hat{\mathbf{z}}$$

$$\nabla \cdot \mathbf{F} = \frac{1}{\rho} \frac{\partial(\rho F_\rho)}{\partial \rho} + \frac{1}{\rho} \frac{\partial F_\phi}{\partial \phi} + \frac{\partial F_z}{\partial z}$$

$$\nabla \times \mathbf{F} = \left(\frac{1}{\rho} \frac{\partial F_z}{\partial \phi} - \frac{\partial F_\phi}{\partial z} \right) \hat{\boldsymbol{\rho}} + \left(\frac{\partial F_\rho}{\partial z} - \frac{\partial F_z}{\partial \rho} \right) \hat{\boldsymbol{\phi}} + \frac{1}{\rho} \left(\frac{\partial(\rho F_\phi)}{\partial \rho} - \frac{\partial F_\rho}{\partial \phi} \right) \hat{\mathbf{z}}$$

$$\nabla^2 f = \frac{1}{\rho} \frac{\partial}{\partial \rho} \left(\rho \frac{\partial f}{\partial \rho} \right) + \frac{1}{\rho^2} \frac{\partial^2 f}{\partial \phi^2} + \frac{\partial^2 f}{\partial z^2}$$

Spherical Polars: $\mathbf{x} = (r \sin \theta \cos \phi, r \sin \theta \sin \phi, r \cos \theta)$

$$\nabla f = \frac{\partial f}{\partial r} \hat{\mathbf{r}} + \frac{1}{r} \frac{\partial f}{\partial \theta} \hat{\boldsymbol{\theta}} + \frac{1}{r \sin \theta} \frac{\partial f}{\partial \phi} \hat{\boldsymbol{\phi}}$$

$$\nabla \cdot \mathbf{F} = \frac{1}{r^2} \frac{\partial(r^2 F_r)}{\partial r} + \frac{1}{r \sin \theta} \frac{\partial(\sin \theta F_\theta)}{\partial \theta} + \frac{1}{r \sin \theta} \frac{\partial F_\phi}{\partial \phi}$$

$$\nabla \times \mathbf{F} = \frac{1}{r \sin \theta} \left(\frac{\partial(\sin \theta F_\phi)}{\partial \theta} - \frac{\partial F_\theta}{\partial \phi} \right) \hat{\mathbf{r}} + \frac{1}{r} \left(\frac{1}{\sin \theta} \frac{\partial F_r}{\partial \phi} - \frac{\partial(r F_\phi)}{\partial r} \right) \hat{\boldsymbol{\theta}} + \frac{1}{r} \left(\frac{\partial(r F_\theta)}{\partial r} - \frac{\partial F_r}{\partial \theta} \right) \hat{\boldsymbol{\phi}}$$

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