

# PHYSICS 204A: MATH METHODS

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Lecture 1.

## Wednesday, September 25, 2019

Reading assignment: read Ch. 2 and 3 of the course text (Arfken/Weber). This is basic linear algebra and vector analysis.

The purpose of this course is to learn formal aspects of quantum mechanics. We'll focus on doing analysis in Hilbert space. It's a remarkable fact about the natural world that most of our physical world is well-approximated by linear systems.

In the simplest form, we may think of vectors as arrays of numbers,

$$(v_1, v_2, \dots). \quad (1.1)$$

But we can also think of some real function  $f(x) : \mathbb{R} \rightarrow \mathbb{R}$  as a collection of numbers too, just by taking its values at arbitrarily close points.

**Definition 1.2.** A *linear vector space* over a field  $F$ , denoted  $L(F)$ , is a set  $\{|v\rangle\}$  with an addition operation  $+$  such that

- for  $|v\rangle, |u\rangle \in L, |u\rangle + |v\rangle = |w\rangle \in L$  (closure)
- for  $c \in F, c|v\rangle \in L$  (scalar multiplication).

These axioms directly imply that any linear combination of vectors in  $L$  is also in the vector space:

$$c_v |v\rangle + c_u |u\rangle \in L. \quad (1.3)$$

This leads us naturally to the notion of *linear (in)dependence*. Suppose we take some vectors  $|v_k\rangle \in L$  and make a linear combo,

$$\sum_k c_k |v_k\rangle. \quad (1.4)$$

**Definition 1.5.** A set of vectors  $\{|v_k\rangle\}$  is *linearly dependent* if there exists some  $\{c_k\}$  not all zero such that

$$\sum_k c_k |v_k\rangle = |0\rangle, \quad (1.6)$$

and such that  $|0\rangle \notin \{|v_k\rangle\}$ .

We need this last condition because otherwise we could simply take the coefficient of the  $|0\rangle$  vector to be 1 and then arrive at a trivial solution.

**Definition 1.7.** If a set of vectors is not linearly dependent, it is *linearly independent*.

The next question we might ask is as follows: what is the size of the biggest set of linearly independent vectors we can construct for a given vector space?

**Definition 1.8.** The maximum number of linearly independent vectors associated to a given vector space is called the *dimension*.

**Example 1.9.** We may consider an infinitely differentiable ( $C^\infty$ ) function. It has a Taylor expansion

$$f(x) = \sum \frac{f^{(n)}(0)}{n!} x^n, \quad (1.10)$$

which we may think of as an expansion in the basis  $(1, x, x^2, \dots)$ .

So this is a vector space with countably infinite dimension. But we can have uncountably infinite-dimensional spaces too, e.g. the space of Fourier-transformable functions in a basis  $e^{ikx}, k \in \mathbb{R}$ . These factors are not just linearly independent; introducing an appropriate inner product, they are orthogonal.

It follows that for a vector space  $L_D$  of dimension  $D$ , any set with more than  $D$  vectors must be linearly dependent, i.e.  $\exists \bar{c}_v, \bar{c}_k$  not all zero such that

$$\bar{c}_v |v\rangle + \sum_k \bar{c}_k |v_k\rangle = 0. \quad (1.11)$$

Moreover  $\bar{c}_v \neq 0$  or else the original set  $|v_k\rangle$  would be linearly dependent. Hence we can divide through, define  $\hat{c}_k = \bar{c}_k / \bar{c}_v$ , and write

$$|v\rangle = \sum_k \hat{c}_k |v_k\rangle. \quad (1.12)$$

That is, we have *decomposed* a general vector  $|v\rangle$  in terms of its components  $\hat{c}_k$  with respect to a basis  $|v_k\rangle$ .

We might now be interested in adding more structure to our vector space. Consider  $L_D$  with  $|v\rangle, |u\rangle$ .

**Definition 1.13.** We define an *inner product* by

$$\langle v|u\rangle : (|v\rangle, |u\rangle) \rightarrow F \quad (1.14)$$

as a map from the input vectors to the field over which the vector space is defined, with the following properties:

- $\langle u|(\lambda|v_1\rangle + \mu|v_2\rangle) = \lambda\langle u|v_1\rangle + \mu\langle u|v_2\rangle$  (linearity)
- $\langle v|u\rangle = \langle u|v\rangle^*$
- $\langle v|v\rangle \geq 0$ , with equality only for the zero vector (positive semi-definite).

If we choose a basis  $\{|v_k\rangle\}$ , then if our vectors  $|u\rangle, |v\rangle$  have some expansion in this basis then by linearity that

$$\langle u|v\rangle = \sum_k \hat{c}_k^v \langle u|v_k\rangle, \quad (1.15)$$

and we can expand each of these inner products as

$$\langle u|v_k\rangle = (\langle v_k|u\rangle)^* = \sum_n c_n^{u*} \langle v_n|v_k\rangle. \quad (1.16)$$

It follows that we can write a general inner product as

$$\langle u|v\rangle = \sum_{n,k} c_k^v c_n^{u*} \langle v_n|v_k\rangle. \quad (1.17)$$

Moreover if we could choose a nice basis which had a special property of *orthogonality* or better yet *orthonormality*, we could reduce this to a single sum

$$\langle u|v \rangle = \sum_k c_k^u c_k^v \quad (1.18)$$

in terms of the components alone.

Lecture 2.

**Monday, September 30, 2019**

Last time, we wrote a general form for the dot (inner) product,

$$\langle u|v \rangle = \sum_{n,m} c_n^{*u} c_m^v \langle \phi_n|\phi_m \rangle, \quad (2.1)$$

where the  $|\phi\rangle$ s are basis vectors and  $u, v$  have expansions

$$|u\rangle = \sum_n c_n^u |\phi_n\rangle. \quad (2.2)$$

This is some quadratic form (it depends only quadratically on the components). And indeed it would be very nice if we could define  $\langle \phi_n|\phi_m \rangle = \delta_{nm}$ , so that our double-sum collapses to a single sum.

**Orthonormality** Let us suppose we start with a basis  $\{|\phi_n\rangle\}$  for a vector space  $L_D$ . We shall show that we can construct a new basis  $\{|\chi_n\rangle\}^D$  such that  $\langle \chi_m|\chi_n \rangle = \delta_{mn}$  has the desired property.

WLOG let us number the basis vectors  $|\phi_1\rangle, |\phi_2\rangle, \dots$  and consider some inner products. The inner product

$$\langle \phi_1|\phi_1 \rangle = N_1 \quad (2.3)$$

is some value  $N_1$ . If  $N_1 = 1$  then we are done; otherwise, define

$$|\chi_1\rangle \equiv \frac{1}{\sqrt{N_1}} |\phi_1\rangle \quad (2.4)$$

so that

$$\langle \chi_1|\chi_1 \rangle = \frac{\langle \phi_1|\phi_1 \rangle}{N_1} = 1. \quad (2.5)$$

Hence  $|\chi_1\rangle$  is a unit vector.

Consider the next vector  $|\phi_2\rangle$ . If

$$\langle \chi_1|\phi_2 \rangle = 0, \quad (2.6)$$

then we can normalize and get

$$|\chi_2\rangle = \frac{|\phi_2\rangle}{\sqrt{N_2}}, \quad (2.7)$$

where  $N_2 = \langle \phi_2|\phi_2 \rangle$ . Otherwise, we first subtract off the projection of the first normalized vector,

$$|\hat{\chi}_2\rangle = |\phi_2\rangle - \langle \chi_1|\phi_2 \rangle |\chi_1\rangle, \quad (2.8)$$

so that

$$\langle \chi_1|\hat{\chi}_2 \rangle = \langle \chi_1|\phi_2 \rangle - \langle \chi_1|\phi_2 \rangle \underbrace{\langle \chi_1|\chi_1 \rangle}_{=1} = 0. \quad (2.9)$$

Hence by our definition,  $|\chi_1\rangle$  and  $|\hat{\chi}_2\rangle$  are orthogonal and we can just normalize. Defining

$$|\hat{\chi}_2\rangle = \langle \hat{\chi}_2|\hat{\chi}_2 \rangle, \quad (2.10)$$

we have

$$|\chi_2\rangle = \frac{|\hat{\chi}_2\rangle}{\sqrt{N_2}}, \quad (2.11)$$

which is a unit vector and normal to  $|\chi_1\rangle$ .

We continue by induction, subtracting off projections and normalizing. This is the *Gram-Schmidt procedure*.<sup>1</sup> Notice also that because  $\langle \phi_n|\phi_m \rangle = \langle \phi_m|\phi_n \rangle^*$ , we can consider values of  $m, n$  to be entries in a

<sup>1</sup>Note that the procedure is a little more subtle in the infinite-dimensional case.

Hermitian matrix. Recalling that  $\langle u|u \rangle \geq 0$ , the norm is perfectly well-defined and indeed we can see that orthonormalization is equivalent to diagonalizing a Hermitian matrix.

**More inner products** On a function space, we can define an inner product

$$\langle f|g \rangle = \int_{-\infty}^{\infty} dx f^*(x)g(x). \quad (2.12)$$

These inner products come with strings attached; our functions usually have to satisfy some integrability properties in order for the inner products to be well-defined. Often the functions we're interested in come from differential equations. Most of the ones we encounter in physics are second-order so these functions ought to be twice-differentiable.<sup>2</sup> And we should also require that our functions are square-integrable so that the integral is well-defined.

We can then define

$$\begin{aligned} \langle u|u \rangle &= \sum_{n,m} c_n^* c_m \langle \phi_n | \phi_m \rangle \\ &= \sum_n |c_n|^2. \end{aligned} \quad (2.13)$$

And this is none other than the generalization of Pythagoras's theorem.

**Schwarz inequality** From the axioms, we can prove the following inequality.

$$|\langle f|g \rangle|^2 \leq \langle f|f \rangle \langle g|g \rangle. \quad (2.14)$$

We may define the linear combination

$$|f - \lambda g \rangle \quad (2.15)$$

and consider its norm (squared)

$$\langle f - \lambda g | f - \lambda g \rangle = \langle f|f \rangle - \lambda^* \langle g|f \rangle - \lambda \langle f|g \rangle + \lambda \lambda^* \langle g|g \rangle. \quad (2.16)$$

This is obviously non-negative, given the axioms. We can extremize this by taking derivatives with respect to  $\lambda, \lambda^*$  (which we may treat as linearly independent, since they are complex) and find that

$$\lambda^* = \frac{\langle g|f \rangle}{\langle g|g \rangle}, \quad \lambda = \frac{\langle f|g \rangle}{\langle g|g \rangle}. \quad (2.17)$$

A bit of manipulation yields the Schwarz inequality.

Let us also note that we can in general translate between the ket notation and vector notation. For a ket vector  $|u \rangle$  we can associate the bra (row) vector  $\langle u| = (c_1^*, c_2^*, \dots, c_D^*)$ . Then the vector inner product is the same as old-fashioned row-column multiplication.

**Bessel inequality** Suppose we rewrite the vector  $|u \rangle$  in a weird way, as

$$|u \rangle = \sum_n' c_n^u |\phi_n \rangle + |\Delta u \rangle, \quad (2.18)$$

where we take some terms and separate them out (so the sum  $\sum'$  omits some indices). We know that

$$|\Delta u \rangle = u - \sum_n' c_n^u |\phi_n \rangle \neq 0, \quad (2.19)$$

so that

$$0 < \langle \Delta u | \Delta u \rangle \quad (2.20)$$

$$= \langle u - \sum_n' c_n^u \phi_n | u - \sum_n' c_n^u \phi_n \rangle \quad (2.21)$$

$$= \langle u|u \rangle - \sum_n' |c_n^u|^2. \quad (2.22)$$

Check the cross-terms with the definition of  $u$  to get this final term. Rearranging, we get the Bessel inequality, which says that

$$\langle u|u \rangle > \sum_n' |c_n^u|^2, \quad (2.23)$$

i.e. the norm of a vector is greater than the partial sums of the squares of the components.

<sup>2</sup>This is a little too strong, actually. They can have finitely many discontinuities and this is still okay.

**Linear operators** We are primarily interested in linear operators, i.e. linear maps from the vector space to itself obeying

$$A(\lambda|\phi\rangle + \mu|\chi\rangle) = \lambda A|\phi\rangle + \mu A|\chi\rangle. \quad (2.24)$$

We can define two operators to be equal if they have the same action on all vectors, i.e.

$$A|\phi\rangle = B|\phi\rangle \quad (2.25)$$

for all  $\phi \in L$ .

In particular there's a nice way that we can rewrite the identity operator, as

$$\mathbb{I} = \sum_n |\phi_n\rangle\langle\phi_n|. \quad (2.26)$$

Let's prove this: by definition,  $\mathbb{I}|u\rangle = |u\rangle$ . On the other side, we see that

$$\sum_n |\phi_n\rangle\langle u| = \sum_n c_n^u |\phi_n\rangle \equiv |u\rangle. \quad (2.27)$$

Provided that  $\{|\phi_n\rangle\}$  is a complete basis, this operator is indeed the identity.

**The delta function** The Dirac delta function is defined in such a way that

$$\int_a^b f(t)\delta(x-t)dt = f(x), \quad (2.28)$$

provided that  $x$  is in the interval  $(a, b)$ . We could think of this as an inner product, however. We have a function and its shadow on the delta function picks out a value. The delta function isn't properly square-integrable, but we may consider it as having a good inner product with functions in our function space (as the limit of some sequence of square-integrable functions, if you like).

Now we'll do something strange. Let us express the delta function in a function basis,

$$\delta(x-t) = \sum_n c_n(t)\phi_n(x). \quad (2.29)$$

The  $t$  dependence must be in the coefficients since the functions themselves are just given. How do we find the coefficients? Just take the integral

$$\int dx \phi_m^*(x)\delta(x-t) = \int dx \sum_n c_n(t)\phi_n(x)\phi_m^*. \quad (2.30)$$

This is super easy to evaluate. On the RHS we have a Kronecker delta  $\delta_{nm}$  by the orthonormality of the basis, and on the left side we have the evaluation of the basis vector  $\phi_m^*$  at  $t$ , i.e.

$$c_m(t) = \phi_m^*(t). \quad (2.31)$$

Hence

$$\delta(x-t) = \sum_n \phi_n^*(t)\phi_n(x). \quad (2.32)$$

We can see that this had to be the case by substituting our expression for the delta function into an integral:

$$f(x) = \int dt \delta(x-t)f(t) = \sum_n \phi_n(x) \int dt \phi_n^*(t)f(t) \quad (2.33)$$

$$= \sum_n \langle\phi_n|f\rangle\phi_n(x), \quad (2.34)$$

which is none other than the components of  $f$  in the basis  $\phi_n$ .

To make our discussion more concrete, let us consider analytic functions which have Taylor expansion

$$f(x) = \sum \frac{f^{(n)}(0)}{n!} x^n, \quad (2.35)$$

defined over the interval  $[-1, 1]$ . Hence  $\{1, x, x^2, x^3, \dots\}$  form a complete basis set for arbitrarily differentiable functions. They are certainly not orthogonal in general, e.g.  $\int_{-1}^1 dx 1 \cdot x^2 \neq 0$ . But we can make them orthonormal with Gram-Schmidt.

Under this inner product, we have

$$\int_{-1}^1 dx 1 \cdot 1 = 2, \quad (2.36)$$

so our first normalized vector is  $1/\sqrt{2}$ . We can check  $x$ :

$$\int_{-1}^1 dx x \cdot x = 2/3, \quad (2.37)$$

so the next normalized vector is  $\sqrt{3/2}x$ . Continuing this way, we see that  $x^2$  and  $x$  are already orthogonal but

$$\int_{-1}^1 dx \frac{1}{\sqrt{2}} x^2 = \frac{2}{3\sqrt{2}}, \quad (2.38)$$

so our first unit vector is not orthogonal to  $x^2$ . We can instead define

$$\hat{x}^2 = x^2 - \sqrt{\frac{2}{3}}. \quad (2.39)$$

which is now orthogonal to the first unit vector  $1/\sqrt{2}$  and to the second unit vector  $\sqrt{3/2}x$ . We can normalize  $\hat{x}^2$  and determine the third unit vector in this set, which is  $\frac{3x^2}{2} - 1$  (we think).

Let us remark that the most general operators that can be diagonalized are *normal* operators, i.e. those satisfying

$$[A, A^\dagger] = 0. \quad (2.40)$$

Clearly, one set of operators that are not normal are the raising and lowering operators, whose commutator is  $[a, a^\dagger] = 1$ .

Lecture 3.

**Wednesday, October 2, 2019**

Today we'll continue discussing operators. We've discussed the identity operator,

$$\mathbb{I} = \sum |\phi_i\rangle \langle \phi_i|, \quad (3.1)$$

which maps any vector into itself. More generally, we can define an operator as follows:

**Definition 3.2.** An *operator* is a map  $A : |\psi\rangle \rightarrow |\bar{\psi}\rangle$  where  $|\psi\rangle, |\bar{\psi}\rangle \in L_D$  are in the same vector space  $L_D$ .<sup>3</sup> A *linear operator* is an operator obeying the linearity property

$$A(\mu|\psi\rangle + \nu|\chi\rangle) = \mu A|\psi\rangle + \nu A|\chi\rangle. \quad (3.3)$$

If we have a set of linear operators, we may define an addition operation on operators as

$$(A + B)|\psi\rangle = A|\psi\rangle + B|\psi\rangle \quad (3.4)$$

and scalar multiplication as

$$(kA)|\psi\rangle = k(A|\psi\rangle). \quad (3.5)$$

Hence we can take linear combinations of linear operators as

$$kA + lB, \quad (3.6)$$

and under this definition we see that linear operators form a vector space.

But there's another way to combine operators, namely by *composition*. That is, given operators  $A$  and  $B$  we can define a new operator  $AB$  defined by the composition

$$AB|\psi\rangle = A(B|\psi\rangle). \quad (3.7)$$

Composition must satisfy certain properties with respect to the other operations we've defined, namely distributivity with respect to addition:

$$A(B + C) = AB + AC. \quad (3.8)$$

Let us note that the product of operators (composition) is generally not commutative,

$$AB \neq BA \quad (3.9)$$

in general. We know this from matrix multiplication.

<sup>3</sup>The math-inclined among us may talk about the space of inner automorphisms on the vector space.

Alternately, we could define a composition rule using a commutator (bracket),

$$[A, B] = AB - BA. \quad (3.10)$$

It's not hard to check that this rule also satisfies distributivity over addition. If we wished, we could also prove (by crunching through the commutators) the Jacobi identity,

$$[[A, B], C] + [[B, C], A] + [[C, A], B] = 0. \quad (3.11)$$

This is related to the Bianchi identity in differential geometry.

Under our standard composition rule, we can define *inverses*. That is, if an operator  $A$  acts as

$$|\psi'\rangle = A|\psi\rangle, \quad (3.12)$$

then the inverse  $A^{-1}$  (if it exists) is the operator such that

$$A^{-1}|\psi'\rangle = |\psi\rangle. \quad (3.13)$$

That is,  $A^{-1}A|\psi\rangle = |\psi\rangle$ .

As we know, not every operator is invertible. Consider the operator which just sends some vector to zero (i.e. it has a null eigenvector). If

$$A|\psi\rangle = 0 \quad (3.14)$$

then the inverse is not well-defined:  $A^{-1}|0\rangle = ?$ .

**Definition 3.15.** The adjoint  $A^\dagger$  of an operator  $A$  is defined by

$$\langle\chi|A^\dagger|\psi\rangle = \langle\psi|A|\chi\rangle^*. \quad (3.16)$$

**Definition 3.17.** If an operator is self-adjoint,  $H^\dagger = H$ , then we call it Hermitian.

Notice that

$$\langle\psi|H|\psi\rangle = \langle\psi|H^\dagger|\psi\rangle = \langle\psi|H|\psi\rangle^* \quad (3.18)$$

by the definition of the adjoint and hermiticity. Therefore the diagonal matrix elements of  $H$  are *real numbers*, i.e. their eigenvalues are real.

We could have also taken an operator which was *anti-Hermitian*,  $H^\dagger = -H$ , which implies that the diagonal elements are instead purely imaginary by the same argument.

**Definition 3.19.** A *unitary* operator is an operator obeying the property

$$U^{-1} = U^\dagger. \quad (3.20)$$

It's clear that we can restrict to the real case, in which case Hermitian matrices become symmetric matrices and anti-Hermitian matrices become antisymmetric (sometimes called skew-symmetric). Our unitary matrices reduce to orthogonal matrices.

**Example 3.21.** Consider the Hilbert space of smooth square-integrable functions over the real line,  $f \in L$ . Our inner product is the integral

$$\int dx f^*(x)g(x). \quad (3.22)$$

Define the operator  $D = -\frac{d}{dx}$ . What is the adjoint  $D^\dagger$ ? We have

$$\int f^*(-\frac{d}{dx})g = f^*g|_a^b + \int (\frac{d}{dx}f^*)g. \quad (3.23)$$

The boundary term vanishes based on the boundary conditions, i.e. given that  $f, g$  vanish at infinity. Taking the complex conjugate to get the adjoint, what is left is

$$\left[ \int (\frac{d}{dx}f^*)g \right]^* = - \int g^*Df. \quad (3.24)$$

So  $D$  is not Hermitian but  $iD$  is (it adds a minus sign to fix the sign in the integration by parts).

Note also that when the integration region is finite, the boundary conditions become nontrivial. However, if  $f$  and  $g$  vanish at the boundary (e.g.  $[-1, 1]$ ) then we restore hermiticity.

Recall we said that we could assign matrix elements to an operator with respect to some set of vectors,

$$\langle \chi | A | \psi \rangle. \quad (3.25)$$

In fact, it's a Sisyphean task to do this for all sets of vectors, but fortunately (thanks to linearity) it suffices to compute the matrix elements in some (complete) basis. With a basis  $|\psi_i\rangle$  we can define

$$\langle \psi_n | A | \psi_k \rangle = A_{nk}. \quad (3.26)$$

For recall that the identity can be written as  $\mathbb{I} = \sum_k |\psi_k\rangle \langle \psi_k|$ , and suppose  $|\psi\rangle$  has some decomposition in the basis

$$|\psi\rangle = \sum_k \langle \psi_k | \psi \rangle |\psi_k\rangle \quad (3.27)$$

Then

$$\begin{aligned} A|\psi\rangle &= \mathbb{I}A\mathbb{I}|\psi\rangle \\ &= \sum_{k,n} |\psi_k\rangle \langle \psi_k | A | \psi_n \rangle \langle \psi_n | \psi \rangle. \end{aligned}$$

Thus we can recognize the components of  $|\psi\rangle$  in our basis, which are given by  $\langle \psi_n | \psi \rangle = C_n$ . Hence

$$\begin{aligned} A \left( \sum_k \langle \psi_k | \psi \rangle |\psi_k\rangle \right) &= \sum_k |\psi_k\rangle \left( \sum_n \langle \psi_k | A | \psi_n \rangle \langle \psi_n | \psi \rangle \right) \\ &= \sum_k |\psi_k\rangle \left( \sum_n A_{kn} C_n \right) \\ &= \sum_k |\psi_k\rangle C_k^{\tilde{\psi}} \end{aligned}$$

in terms of the components of some new vector  $|\tilde{\psi}\rangle$ . That is, if we know the matrix elements of  $A$  in some basis and we know the components of the vector in that basis, we can uniquely determine the components of its image under  $A$  in the same basis.

In the end, this is just abstract matrix multiplication. That is,

$$C_n^{\tilde{\psi}} = \sum_k A_{nk} C_k^{\psi}. \quad (3.28)$$

We can also write the operator  $A$  in terms of its matrix elements:

$$\begin{aligned} A &= \mathbb{I}A\mathbb{I} \\ &= \sum_{n,k} |\psi_k\rangle \langle \psi_k | A | \psi_n \rangle \langle \psi_n | \\ &= \sum_{n,k} A_{kn} |\psi_k\rangle \langle \psi_n|. \end{aligned}$$

This also tells us immediately that the matrix elements of the identity in any orthonormal basis are as we could have guessed—  $\mathbb{I}_{nk} = \delta_{nk}$ , the Kronecker delta.

It also follows that the matrix elements of the adjoint of an operator obey

$$(A^\dagger)_{nk} = A_{kn}^*. \quad (3.29)$$

This gives us another statement of hermiticity— equivalently, a hermitian operator is one whose matrix elements obey

$$A_{kn}^* = A_{nk}. \quad (3.30)$$

And thus

$$A^\dagger = \sum_{k,n} A_{nk}^* |\psi_k\rangle \langle \psi_n|. \quad (3.31)$$

Let's check that for hermitian operators, the expectation value is non-negative,

$$\langle \psi | A | \psi \rangle \geq 0. \quad (3.32)$$



Writing  $A$  in terms of its matrix elements, we have

$$\begin{aligned}\langle\psi|A|\psi\rangle &= \sum_{n,k} \langle\psi|\psi_k\rangle A_{kn} \langle\psi_n|\psi\rangle \\ &= C_k^{*\psi} A_{kn} C_n^\psi.\end{aligned}$$

Suppose we have two orthonormal bases for the same space,  $\{|\psi_k\rangle\}, \{|\psi'_k\rangle\}$ . It follows that the new basis has some decomposition in the old basis. That is, the set  $\{|\psi_k\rangle, |\psi'_1\rangle\}$  is linearly dependent and so

$$|\psi'_n\rangle = \sum_k c_k^{n'} |\psi_k\rangle, \quad (3.33)$$

in terms of some coefficients  $c_k^{n'}$ . It's also true that we can go back,

$$|\psi_n\rangle = \sum_k c_k^n |\psi'_k\rangle. \quad (3.34)$$

Certainly we can write this decomposition as the action of an operator  $U$ :

$$U|\psi_n\rangle = \sum_k c_k^{n'} |\psi_k\rangle, \quad (3.35)$$

and moreover  $U$  must be invertible.

Lecture 4.

**Monday, October 7, 2019**

When we choose coordinates, the name of the game is to exploit the *symmetries* of the problem. That is, to find coordinates which respect the dynamical symmetries of the Hamiltonian. If we choose our coordinates well enough, the equations of motion become trivial to solve.

Suppose we have two bases  $\{|\psi_i\rangle\}, \{|\psi'_i\rangle\}$ . Then we can certainly write

$$U|\psi_i\rangle = |\psi'_i\rangle = \sum_j C_{ij}^{\psi'} |\psi_j\rangle \quad (4.1)$$

or equivalently

$$\bar{U}|\psi'_j\rangle = |\psi_j\rangle = \sum_i C_{jk}^{\psi} |\psi'_i\rangle. \quad (4.2)$$

That is, we can express an element of one basis in another basis. It is evident that the inverse exists, since it doesn't matter what we call the first and the second basis. So take the first equation and act on it with  $U^{-1}$ . Then

$$|\psi_i\rangle = U^{-1}U|\psi_i\rangle = U^{-1}|\psi'_i\rangle = \sum_j C_{ij}^{\psi'} U^{-1}|\psi_j\rangle. \quad (4.3)$$

What are the matrix elements  $\langle\psi_k|U|\psi_i\rangle = \langle\psi_k|\psi'_i\rangle$ ?

$$\langle\psi_k|\psi'_i\rangle = \sum_j C_{ij}^{\psi'} \langle\psi_k|\psi_j\rangle, \quad (4.4)$$

or equivalently

$$|\psi'_i\rangle = \sum_j |\psi_j\rangle \langle\psi_j|\psi'_i\rangle. \quad (4.5)$$

Similarly

$$|\psi_j\rangle = \sum_i |\psi'_i\rangle \langle\psi'_i|\psi_j\rangle. \quad (4.6)$$

Notice also that these coefficients are therefore clearly related by

$$\langle\psi'_i|\psi_j\rangle = \langle\psi_j|\psi'_i\rangle^*. \quad (4.7)$$

- Consider now the inner product

$$\begin{aligned}
 \delta_{kj} &= \langle \psi_k | \psi_j \rangle \\
 &= \sum_{i,l} \langle \psi'_l | \psi_k \rangle^* \underbrace{\langle \psi'_l | \psi'_i \rangle}_{\delta_{li}} \langle \psi'_i | \psi_j \rangle \\
 &= \sum_i \langle \psi'_i | \psi_k \rangle^* \langle \psi'_i | \psi_j \rangle \\
 &= \sum_i \langle \psi_k | \psi'_i \rangle \langle \psi'_i | \psi_j \rangle.
 \end{aligned}$$

Define

$$U_{ij} = \langle \psi'_i | \psi_j \rangle. \quad (4.8)$$

And similarly

$$\bar{U}_{ki} = \langle \psi_k | \psi'_i \rangle = U_{ik}^*. \quad (4.9)$$

Hence

$$\delta_{kj} = \sum_i U_{ki}' U_{ij}, \quad (4.10)$$

so since  $(U^{-1})_{kj} = U_{jk}^*$ , we see that transformations between two orthonormal bases are unitary, i.e. they satisfy

$$UU^\dagger = \mathbb{I}. \quad (4.11)$$

We could have seen this in a basis-free way by requiring that inner products do not depend on the choice of basis. Hence

$$\langle \chi' | \psi' \rangle = \langle \chi | U^\dagger U | \psi \rangle = \langle \chi | \psi \rangle = \langle \chi | \mathbb{I} | \psi \rangle, \quad (4.12)$$

and hence it must be that  $U^\dagger U = \mathbb{I}$  since they agree on *any* vectors  $|\chi\rangle, |\psi\rangle$ .

We've concluded that changes of basis can be written as unitary transformations and moreover that inner products must be invariant under such transformations. What about operators? Consider

$$A = \sum_{i,j} a_{ij} |\psi_i\rangle \langle \psi_j|. \quad (4.13)$$

Suppose we have a new vector  $|\phi\rangle$  given by

$$|\phi\rangle = \sum_i |\psi_i\rangle \langle \psi_i | \phi \rangle, \quad (4.14)$$

and we act on it with a unitary  $U$  to get some new  $|\phi'\rangle$ . Hence

$$U|\phi\rangle = \sum_i U|\psi_i\rangle \langle \psi_i | \phi \rangle. \quad (4.15)$$

Let  $U$  take us between bases, such that  $|\psi'_i\rangle = U|\psi_i\rangle$  or equivalently

$$U^\dagger |\psi'_i\rangle = |\psi_i\rangle. \quad (4.16)$$

The easy way to see what happens to operators is to recognize that

$$U|\phi'\rangle = UA|\phi\rangle = UAU^\dagger(U|\phi\rangle), \quad (4.17)$$

so that

$$A' = UAU^\dagger \quad (4.18)$$

is the corresponding operator to  $A$  in the new basis. Looking at the spectral decomposition,

$$A' = UAU^\dagger = \sum U|\psi_i\rangle A_{ij} \langle \psi_j | U^\dagger \quad (4.19)$$

$$= \sum |\psi'_i\rangle A_{ij} \langle \psi'_j|, \quad (4.20)$$

so the matrix elements are left unchanged in the new basis. In fact, it must have been so, since

$$\langle \psi_i | A | \psi_j \rangle = \left( \langle \psi_i | U^\dagger \right) \left( UAU^\dagger \right) (U|\psi_j\rangle). \quad (4.21)$$

Let us now make a connection to quantum mechanics. If the Hamiltonian is time-independent, recall that we can write the solution of

$$i\partial_t|\psi\rangle = H|\psi\rangle \quad (4.22)$$

as

$$|\psi(t)\rangle = e^{-itH}|\psi(0)\rangle. \quad (4.23)$$

That is, we may think of time evolution as a complex rotation of the initial condition.<sup>4</sup>

Consider now a set of vectors which are not orthonormal but are linearly independent,  $\{|\chi_i\rangle\}$ . By Gram-Schmidt we can therefore define  $\{|\psi_\mu\rangle\}$  as the result of acting on  $|\psi_i\rangle$  with an operator  $T$  such that

$$|\psi_\mu\rangle = \sum_{i=1}^{\mu} T_{\mu i} |\chi_i\rangle. \quad (4.24)$$

In particular we may take  $T_{\mu i}$  to be upper-triangular because we limit the sum to run only up to  $\mu$ .

Now observe that

$$\begin{aligned} \delta_{\nu\mu} &= \langle\psi_\nu|\psi_\mu\rangle \\ &= \sum_{ij} T_{\nu j}^* T_{\mu i} \langle\chi_j|\chi_i\rangle \\ &= \sum_{ij} T_{\nu j}^* \langle\chi_j|\chi_i\rangle T_{\mu i} \\ &= (T^\dagger S T)_{\nu\mu}. \end{aligned}$$

These final inner products are in general not 1; they are like a metric, in that they relate the original basis vectors. We see that

$$\mathbb{I} = T^\dagger S T. \quad (4.25)$$

But notice that  $T$  is invertible, and therefore so is  $T^\dagger$ . We find that

$$(T^\dagger)^{-1} T^{-1} = S, \quad (4.26)$$

so the matrix  $S$  relating the original vectors can be written in terms of the matrix  $T$  which performs the orthonormalization process. In general, this allows us to perform a *similarity transformation* so that we can write some operator  $A' = T A T^{-1}$  and

$$A = (T^\dagger)^{-1} \bar{A} T^{-1}, \quad (4.27)$$

which says that an operator may be written as the product of an upper triangular matrix, a diagonal matrix, and a lower triangular matrix. This is a special case of the Jordan decomposition of an operator.

Notice also that unitaries leave the trace and the determinant unchanged:

$$\text{Tr}(U A U^\dagger) = \text{Tr}(A U^\dagger U) = \text{Tr}(A), \quad (4.28)$$

and determinants are unchanged since

$$U U^\dagger = \mathbb{I} \implies 1 = \det \mathbb{I} = \det U \det U^\dagger = (\det U)(\det U)^*, \quad (4.29)$$

so the determinant of  $U$  is a phase  $e^{i\delta}$  and

$$\det(U A U^\dagger) = \det(A) \det(U^\dagger U) = \det A. \quad (4.30)$$

This helps our intuition that unitary matrices are really complex generalizations of rotations (orthogonal matrices).

It will often be our interest in quantum mechanics of solving the eigenvector and eigenvalue problem, i.e.

$$A|\phi\rangle = \lambda|\phi\rangle. \quad (4.31)$$

Thanks to the fundamental theorem of algebra, we are guaranteed at least one eigenvalue, some  $\lambda$  satisfying the characteristic equation.

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<sup>4</sup>There is a book describing the structure of path integrals in quantum mechanics, by Feynman and Hibbs, related to the exponentiation formula for the Hamiltonian. There is also a text by Messiah which contains lots of good quantum mechanics.

Lecture 5.

**Wednesday, October 9, 2019**

*"If you can ask a general-sounding question that will help you with the homework, by all means do it. I encourage you to be sneaky."* –Nemanja Kaloper

We said last time that operators in a new basis  $|\phi'_i\rangle = U|\phi_i\rangle$  are given by

$$A' = UAU^\dagger = \sum a_{ik}|\phi'_i\rangle\langle\phi'_k|. \quad (5.1)$$

Hence the *components* of the corresponding operator  $A'$  are left unchanged in the new basis. However, it is also true that the original operator has some components<sup>5</sup>

$$A = \sum a_{ik}|\phi_i\rangle\langle\phi_k| = \sum a'_{ik}|\phi'_i\rangle\langle\phi'_k|. \quad (5.2)$$

That is, we can either define a change of basis and also rotate the operator along with it ( $A' = UAU^\dagger$ ) or we can just change the basis and leave the operator fixed. Hence its projections onto the new basis will change and therefore its components will change. This is the difference between *active and passive transformations*. Incidentally, this is related to the Schrödinger and Heisenberg pictures. There is also a third hybrid<sup>6</sup> picture known as the interaction picture, and this ends up being the most practical one to do perturbative calculations in.

We will work mostly in the picture where operators are fixed and the bases change. How do we find the components in the new basis?

$$|\phi'_i\rangle = U|\phi_i\rangle \quad (5.3)$$

$$= \sum_j |\phi_j\rangle\langle\phi_j|U|\phi_i\rangle \quad (5.4)$$

$$= \sum_j U_{ji}|\phi_j\rangle. \quad (5.5)$$

Hence

$$a'_{ik} = \sum_{jl} \langle\phi'_i|\phi_j\rangle a_{jl} \langle\phi_l|\phi'_k\rangle = \sum_{jl} U_{ji}^* a_{jl} U_{lk}, \quad (5.6)$$

which tells us that the components transform in the opposite way from the operator. That is, in this picture we have

$$A_{ik}^{\phi'} = a'_{ik} = (U^\dagger A U)_{ik}. \quad (5.7)$$

Recall from last time that for some general operator  $A$ , we are guaranteed at least one eigenvector,

$$A|\psi\rangle = \lambda|\psi\rangle. \quad (5.8)$$

Suppose we have a coupled harmonic oscillator potential

$$V = ax^2 + bxy + cy^2, \quad (5.9)$$

which results in the forces

$$F_x = -2ax - by \quad (5.10)$$

$$F_y = -2cy - bx. \quad (5.11)$$

This looks like an ellipsoidal well, and particles will form closed orbits (Lissajous figures) in this well.

In general orbits will be 2-dimensional, changing in  $x$  and  $y$ . But! Sometimes if we set it up right, we will find that our orbits become 1-dimensional. And this tells us that we've found the *characteristic directions* of the oscillator.

We can now write this as an eigenvalue problem:

$$\begin{pmatrix} a_x \\ a_y \end{pmatrix} = \frac{1}{m} \begin{pmatrix} F_x \\ F_y \end{pmatrix} = \begin{pmatrix} -2ax - by \\ -2cy - bx \end{pmatrix} = \begin{pmatrix} -2a & -b \\ -b & -2c \end{pmatrix} \begin{pmatrix} x \\ y \end{pmatrix}. \quad (5.12)$$

<sup>5</sup>I tend to write this as components times  $|\phi_i\rangle\langle\phi_k|$  to emphasize that this is really a linear combination of operators.

<sup>6</sup>"Bastard," as per Nemanja.

We could imagine an orbit where the acceleration becomes parallel to the displacement, in which case the motion just becomes one-dimensional.

If we project down the equipotentials, we can read off the principal axes, which tells us the right directions in which the motion separates.

We can abstract the problem: consider

$$\begin{pmatrix} h_{11} & h_{12} \\ h_{12} & h_{22} \end{pmatrix} \begin{pmatrix} x \\ y \end{pmatrix} = \lambda \begin{pmatrix} x \\ y \end{pmatrix}. \quad (5.13)$$

We solve the eigenvalue problem by finding the characteristic equation. That is,

$$(h_{11} - \lambda)(h_{22} - \lambda) - h_{12}^2 = 0. \quad (5.14)$$

These could be real eigenvalues, and potentially degenerate.<sup>7</sup> In this example, the matrix is real and symmetric and therefore Hermitian, which implies its eigenvalues are real. In the general case the eigenvalues could be complex if the matrix entries were not all real.<sup>8</sup> Note that we get the constraint equations,

$$(h_{11} - \lambda_1)x + h_{12}y = 0 \quad (5.15)$$

$$h_{12}x + (h_{22} - \lambda_1)y = 0. \quad (5.16)$$

Notice that the determinant of  $H - \lambda \mathbb{I}$  vanished, which tells us that the rows are linearly *dependent*. So we will not get any extra information out of the second equation, i.e. our solution is not uniquely determined.<sup>9</sup> We can solve

$$y = -\frac{h_{11} - \lambda_1}{h_{12}}x, \quad (5.17)$$

but we cannot a priori fix the value of  $x$ . For notice that to any eigenvector, i.e. a solution to

$$\begin{pmatrix} h_{11} - \lambda_1 & h_{12} \\ h_{12} & h_{22} - \lambda_1 \end{pmatrix} \begin{pmatrix} x \\ y \end{pmatrix} = 0, \quad (5.18)$$

another (nontrivial) solution to this is clearly  $\begin{pmatrix} cx \\ cy \end{pmatrix}$  for some  $c \neq 0$ . Hence eigenvectors are only determined up to an overall normalization factor.

For instance, let us compute the eigenvectors of

$$\begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix}. \quad (5.19)$$

The characteristic equation is

$$\lambda^2 - 1 \implies \lambda = \pm 1. \quad (5.20)$$

Hence we have

$$\begin{pmatrix} y \\ x \end{pmatrix} = \pm \begin{pmatrix} x \\ y \end{pmatrix}, \quad (5.21)$$

giving  $y = x$  or  $y = -x$ . The normalized eigenvectors are

$$\frac{1}{\sqrt{2}} \begin{pmatrix} 1 \\ 1 \end{pmatrix}, \frac{1}{\sqrt{2}} \begin{pmatrix} 1 \\ -1 \end{pmatrix}. \quad (5.22)$$

Moreover, notice that we can build a matrix built out of the components of the eigenvectors, namely

$$E = \begin{pmatrix} x_+ & x_- \\ y_+ & y_- \end{pmatrix} = \frac{1}{\sqrt{2}} \begin{pmatrix} 1 & 1 \\ 1 & -1 \end{pmatrix}. \quad (5.23)$$

Observe that

$$\begin{pmatrix} 0 & 1 \\ 1 & 0 \end{pmatrix} E = E \begin{pmatrix} +1 & 0 \\ 0 & 1 \end{pmatrix} = E \begin{pmatrix} \lambda_+ & 0 \\ 0 & \lambda_- \end{pmatrix}. \quad (5.24)$$

<sup>7</sup>In this case, if the eigenvalues are degenerate, then our elliptical bowl becomes a circle. There are not just two characteristic directions but a continuum of characteristic directions.

<sup>8</sup>Incidentally they would be complex conjugates, since complex roots to real equations must come in pairs.

<sup>9</sup>In fancier language, we get a one-parameter family of eigenvalues, i.e. a 1D subspace of eigenvalues.

Moreover,  $E^T E = \mathbb{I}$  since the eigenvectors are orthonormal. So if we now multiply on the left by  $E^T$  we find that

$$E^T A E = A' \quad (5.25)$$

where  $A'$  is now the diagonal matrix of the eigenvalues of  $A$ .

Lecture 6.

**Monday, October 14, 2019**

Last time, we showed that certain kinds of matrices can be diagonalized, i.e. we can write them as a similarity transform of a diagonal matrix. The goal is basically to project onto the (normalized) eigenvectors ( $E$ ), multiply by eigenvalues in the eigenbasis ( $D$ ), and then perform the inverse transformation to take us back to the original basis ( $E^T$ ).

Recall we are interested in the eigenvalue problem for Hermitian matrices,

$$H|\psi\rangle = \lambda_\psi |\psi\rangle \quad (6.1)$$

where  $H = H^\dagger$ . In the finite dimensional case, we are guaranteed at least one (simply write down the characteristic equation, and it must have at least one root over the complex numbers). Suppose moreover we have another eigenvalue

$$H|\phi\rangle = \lambda_\phi |\phi\rangle. \quad (6.2)$$

Let us now conjugate the second equation to get

$$\langle\phi|H = \lambda_\phi^* \langle\phi|. \quad (6.3)$$

We can multiply Eqn. 6.1 by  $\langle\phi|$  and Eqn. 6.3 by  $|\psi\rangle$ . Hence

$$\langle\phi|H|\psi\rangle = \lambda_\psi \langle\phi|\psi\rangle \quad (6.4)$$

$$\langle\phi|H|\psi\rangle = \lambda_\phi^* \langle\phi|\psi\rangle. \quad (6.5)$$

We can now subtract these two from each other to find that

$$0 = (\lambda_\psi - \lambda_\phi^*) \langle\phi|\psi\rangle. \quad (6.6)$$

If  $|\phi\rangle = |\psi\rangle$  with  $|\psi\rangle \neq 0$ , then we immediately find that

$$\lambda_\phi = \lambda_\phi^*, \quad (6.7)$$

which tells us the eigenvalues are real.

Conversely if the eigenvalues are distinct ( $\lambda_\psi \neq \lambda_\phi$ ) then

$$0 = (\lambda_\psi - \lambda_\phi) \langle\phi|\psi\rangle \implies \langle\phi|\psi\rangle = 0. \quad (6.8)$$

Hence eigenvectors associated to distinct eigenvalues *must be orthogonal*.

Finally, suppose we have a degenerate case where multiple eigenvectors are associated to the same eigenvalue,  $\lambda_\phi = \lambda_\psi$ . Then this equation cannot help us because it automatically vanishes. We get a degenerate subspace associated to the eigenvalue  $\lambda_\phi$ , and we can then perform Gram-Schmidt within the subspace.

This completes the process of diagonalization. Distinct eigenvalues correspond to orthogonal eigenvectors, and degenerate eigenvalues give us a subspace whose basis vectors can be made orthogonal by Gram-Schmidt.

Moreover, there is a completeness relation for the eigenvectors:

$$\sum_i |\psi_i\rangle \langle\psi_i| = \mathbb{I}. \quad (6.9)$$

This is fairly trivial in the finite-dimensional case (in  $n$  dimensions we can construct  $n$  orthogonal vectors from the eigenvectors, so we get a basis). However, there is a convergence requirement in the infinite-dimensional case. It is not enough to just count and show orthogonality; we must show that all the same information from the original function is contained in the Fourier coefficients, for example.

We will keep in the back of our mind that this completeness relation is really something we have to prove, but for the most part we will perform calculations as though it is true. That is, recall we can write a decomposition in some basis as

$$|\phi\rangle = \sum_i \langle\psi_i|\phi\rangle |\psi_i\rangle. \quad (6.10)$$

If

$$\lim_{N \rightarrow \infty} \left\| |\phi\rangle - \sum_i^N \langle\psi_i|\phi\rangle |\psi_i\rangle \right\|^2 = 0 \quad (6.11)$$

then we may say that our basis is complete in the sense that a decomposition into components  $\langle\psi_i|\phi\rangle$  converges to the “real” vector  $|\phi\rangle$ .

Now

$$H = H \sum_i |\psi_i\rangle \langle\psi_i| \quad (6.12)$$

$$= \sum_i H |\psi_i\rangle \langle\psi_i| \quad (6.13)$$

$$= \sum_i \lambda_i |\psi_i\rangle \langle\psi_i|. \quad (6.14)$$

That is,  $H$  is diagonal in its eigenbasis and its elements are the eigenvalues. Let us note that

$$H \begin{pmatrix} \psi_1^1 & \psi_1^2 \dots \\ \vdots & \vdots \\ \psi_N^1 & \psi_N^2 \dots \end{pmatrix} = \begin{pmatrix} \psi_1^1 & \psi_1^2 \dots \\ \vdots & \vdots \\ \psi_N^1 & \psi_N^2 \dots \end{pmatrix} \begin{pmatrix} \lambda_1 & & \\ & \lambda_2 & \\ & & \ddots \end{pmatrix}, \quad (6.15)$$

where lower indices indicate which eigenvector we’re looking at and upper indices indicate the component of that eigenvector. That is, this is just a matrix form of

$$H|\psi_i\rangle = \lambda_i |\psi_i\rangle. \quad (6.16)$$

More generally we have constructed

$$HU = UH_d \implies H_d = U^\dagger HU, \quad (6.17)$$

which tells us that  $U$  gives the complex rotation to take us to the eigenbasis of the hermitian operator  $H$ .

Now suppose that two operators commute,  $[A, B] = 0$ . Then a mutual eigenbasis can be found:

$$[A, B] \iff A|\phi\rangle = a|\phi\rangle, B|\phi\rangle = b|\phi\rangle, \quad (6.18)$$

for all eigenvectors  $|\phi\rangle$ . The  $\iff$  direction is easy:

$$AB|\phi\rangle = bA|\phi\rangle = ab|\phi\rangle, \quad (6.19)$$

and similarly

$$BA|\phi\rangle = ba|\phi\rangle = ab|\phi\rangle \quad (6.20)$$

for scalars. By completeness,

$$(AB - BA)|\psi\rangle = 0 \quad (6.21)$$

for a general vector  $|\psi\rangle$ .

Let us now prove the  $\implies$  direction. If  $A$  and  $B$  commute, then

$$AB|\phi\rangle = BA|\phi\rangle \quad (6.22)$$

for all  $|\phi\rangle$ . Let us suppose  $|\phi\rangle$  is an eigenvector of  $B$ ,

$$B|\phi\rangle = b|\phi\rangle. \quad (6.23)$$

Then

$$bA|\phi\rangle = BA|\phi\rangle. \quad (6.24)$$

But this tells us that  $A|\phi\rangle$  is an eigenvector of  $B$  with eigenvalue  $b$ .

There are a few possibilities here. It could be that  $A|\phi\rangle = 0$ , which tells us that  $|\phi\rangle$  is an eigenvector of  $A$  with eigenvalue zero,

$$A|\phi\rangle = 0|\phi\rangle. \quad (6.25)$$

The next case is where  $b$  is a unique eigenvalue of  $B$ . Then

$$A|\phi\rangle \propto |\phi\rangle, \quad (6.26)$$

i.e. it lives in the same 1D subspace as  $|\phi\rangle$ . It is simply proportional to the original eigenvector  $|\phi\rangle$ , and hence

$$A|\phi\rangle = a|\phi\rangle. \quad (6.27)$$

The only subtle case is when  $b$  is a degenerate eigenvalue. What then? In this case,  $A|\phi\rangle$  simply lives in the degenerate subspace corresponding to eigenvalue  $b$ .

**Normal operators** Suppose we have an operator  $A$  obeying

$$[A, A^\dagger] = 0. \quad (6.28)$$

Suppose  $A$  has an eigenvalue,

$$A|\phi_1\rangle = \lambda_1|\phi_1\rangle \implies (A - \lambda_1)|\phi_1\rangle = 0. \quad (6.29)$$

Taking the dagger,

$$\langle\phi_1|(A^\dagger - \lambda_1^*) = 0. \quad (6.30)$$

Now since these are individually zero, we can multiply them to get

$$\langle\phi_1|(A^\dagger - \lambda_1^*)(A - \lambda_1)|\phi_1\rangle = 0. \quad (6.31)$$

Expanding out the multiplication we have

$$\begin{aligned} 0 &= \langle\phi_1|(A^\dagger A - \lambda_1^* A - \lambda A^\dagger + \lambda_1^* \lambda_1)|\phi_1\rangle \\ &= \langle\phi_1|(A A^\dagger - \lambda_1^* A - \lambda A^\dagger + \lambda_1^* \lambda_1)|\phi_1\rangle \\ &= \langle\phi_1|(A - \lambda_1)(A^\dagger - \lambda_1^*)|\phi_1\rangle \end{aligned}$$

where in the second line we have used the fact that  $A$  and  $A^\dagger$  commute. This is just the modulus  $\|(A^\dagger - \lambda_1^*)|\phi_1\rangle\|^2$ , so since it is zero,

$$(A^\dagger - \lambda_1^*)|\phi_1\rangle = 0 \implies A^\dagger|\phi_1\rangle = \lambda_1^*|\phi_1\rangle. \quad (6.32)$$

We conclude that  $|\phi_1\rangle$  is an eigenvector of  $A^\dagger$  with eigenvalue  $\lambda_1^*$ .

Consider now some vector  $|\psi\rangle$  in the orthogonal complement of  $|\phi_1\rangle$ ,

$$|\psi\rangle \in L_D \setminus \{|\phi\rangle | |\phi\rangle = a|\phi_1\rangle, a \in \mathbb{C}\}. \quad (6.33)$$

Then

$$0 = \langle\phi_1|\psi\rangle \quad (6.34)$$

by definition, and

$$\langle\phi_1|A|\psi\rangle = \langle\psi|A^\dagger|\phi_1\rangle^* = [\lambda_1^*\langle\psi|\phi_1\rangle]^* = \lambda_1\langle\phi_1|\psi\rangle = 0, \quad (6.35)$$

since  $|\phi_1\rangle$  is an eigenvector of  $A^\dagger$ . Hence it is not just  $|\psi\rangle$  that is in the orthogonal complement of  $|\phi_1\rangle$ —in fact, its image  $A|\psi\rangle$  is also in the orthogonal complement. Hence this orthogonal complement is well-defined. Its image  $A|\psi\rangle$  does not mix with any of the vectors we just took away.

Now we have reduced the dimension of the space by one, which tells us that we can repeat this process with the resulting subspace and therefore find a complete set of eigenvectors and eigenvalues. This completes the proof that normal operators can be diagonalized.<sup>10</sup>

As a practical aside, these sorts of techniques appear in the case of perturbation theory. Sometimes we will have a degenerate eigenvalue of the zeroth order Hamiltonian due to certain symmetries of the problem, and if we are lucky, the perturbation will lift (break) this symmetry, splitting the degenerate energy levels. This is the case with the Stark effect (electric field aligning the electric dipole of an electron) and Zeeman effect (magnetic equivalent).

Now suppose we have an eigenbasis in hand and can write

$$H = \sum |\phi_i\rangle \lambda_i \langle\phi_i|. \quad (6.36)$$

<sup>10</sup>In the continuous case we must also be careful about questions of convergence in the inner products so they are well-defined, but this otherwise totally defines the procedure.



Hence

$$H^n |\phi_i\rangle = \lambda_i^n |\phi_i\rangle, \quad (6.37)$$

so in general

$$H^n = \sum |\phi_i\rangle \lambda_i^n \langle \phi_i| \quad (6.38)$$

and we can define a function of an operator by its Taylor expansion,

$$f(H) = \sum_n \frac{f^{(n)}(0)}{n!} H^n \quad (6.39)$$

$$= \sum_i |\phi_i\rangle \sum_n \frac{f^{(n)}(0)}{n!} \lambda_i^n \langle \phi_i| \quad (6.40)$$

$$= \sum_i |\phi_i\rangle f(\lambda_i) \langle \phi_i|, \quad (6.41)$$

provided that the set is complete (so we can change the order of summation over  $i, n$ ) and the  $\lambda_i$ s are in the radius of convergence of  $f$  (so we can perform the sum over  $n$ ).

We can then define the inverse of a Hermitian operator acting on an eigenvector of  $H$ :

$$H^{-1} |\phi_i\rangle = \frac{1}{\lambda_i} |\phi_i\rangle, \quad (6.42)$$

which is valid so long as *none of the eigenvalues are zero*. Then its action on any vector is well-defined since the  $|\phi_i\rangle$  are complete.

Inserting the identity, we can also find that

$$\langle \phi | H | \phi \rangle = \sum \langle \phi | \phi_i \rangle \lambda_i \langle \phi_i | \phi \rangle = \sum_i \lambda_i |\langle \phi_i | \phi \rangle|^2. \quad (6.43)$$

When our states are normalized, this gives us a probabilistic interpretation of the expectation value  $\langle \phi | H | \phi \rangle$  in terms of the probability  $|\langle \phi_i | \phi \rangle|^2$  of measuring  $\lambda_i$ .

**Example 6.44.** Suppose we have three masses attached on springs, with two small masses  $m$  at the sides and a larger mass  $M$  in the center. The spring constants are  $k$ . Label their positions by  $x_1, x_2, x_3$ . Then the potential is

$$V = \frac{k}{2} [(x_1 - x_2)^2 + (x_2 - x_3)^2], \quad (6.45)$$

and the restoring forces are given by

$$\ddot{x}_1 = -\frac{k}{m}(x_1 - x_2) \quad (6.46)$$

$$\ddot{x}_2 = -\frac{k}{M}(x_2 - x_1) - \frac{k}{M}(x_2 - x_3) \quad (6.47)$$

$$\ddot{x}_3 = -\frac{k}{m}(x_3 - x_2) \quad (6.48)$$

We take the ansatz

$$x_i \sim e^{i\omega t} |\phi_i\rangle, \quad (6.49)$$

so that all the derivatives become  $-\omega^2$ s, and then we can easily write this as a matrix equation. The solution for the normal modes reduces to finding the eigenvalues of the matrix

$$\begin{pmatrix} \frac{k}{m} & -\frac{k}{m} & 0 \\ -\frac{k}{M} & \frac{2k}{M} & -\frac{k}{M} \\ 0 & -\frac{k}{m} & \frac{k}{m} \end{pmatrix}. \quad (6.50)$$

Note this matrix is normal (it commutes with its transpose) but not hermitian. That's the standard way to solve this. But we'll try something different. We have translational symmetry—the forces depend only on relative separations. Hence there is an  $\omega^2 = 0$  mode corresponding to the eigenvector  $(1, 1, 1)$ .

There is also another momentum-conserving mode  $(1, 0, -1)$  corresponding to the outer two masses moving, and we can read off the eigenvalue immediately:  $\omega^2 = k/m \implies \omega = \sqrt{k/m}$ .

Finally, there is a third mode. It is orthogonal to the other two, and it takes more work to find:  $(1, -2m/M, 1)$ , with eigenvalue  $\frac{k}{m} + \frac{2k}{M}$ . Hence we could take a cross product and normalize, or we could use Gram-Schmidt to calculate the final eigenvector.

Lecture 7.

### Wednesday, October 16, 2019

Today we'll look at roughly the material from Arfken 7.1-7.4. The topic of this section will be ordinary differential equations. We basically know that the way to solve differential equations is to either perform some integral transformation to make it easier or alternately to know a special function which solves the equation.

A linear ordinary differential equation takes the form

$$\mathcal{L}y(x) = f(x), \quad (7.1)$$

where  $\mathcal{L}$  is a linear operator,

$$\mathcal{L} = p_0(x) + p_1(x)D + \dots + p_n(x)D^n \quad (7.2)$$

where  $D = \frac{d}{dx}$  and  $f$  is the source function. A solution  $y(x)$  can be written as

$$y(x) = \sum_{i=1}^n c_i y_n^{(i)}(x) + y_p(x), \quad (7.3)$$

where the  $y_n$  are solutions to the homogeneous equation (i.e. they are in the kernel of  $\mathcal{L}$ ,  $\mathcal{L}y_n(x) = 0$ ) and  $y_p$  is the "particular solution" fitting the source. It's a general fact that  $\dim(\ker \mathcal{L}) = n$ , i.e. there are  $n$  linearly independent solutions to the homogeneous equation.

**First-order equations** A first order equation looks like

$$y'(x) = f(x, y) = -\frac{P(x, y)}{Q(x, y)}, \quad (7.4)$$

where  $P$  and  $Q$  are some functions. The simplest case is

$$y' = -\frac{P(x)}{Q(y)}, \quad (7.5)$$

which is separable and has implicit solution

$$\int dy Q(y) = - \int dx P(x). \quad (7.6)$$

But we can also rearrange<sup>11</sup> to get

$$P(x, y)dx + Q(x, y)dy = 0. \quad (7.7)$$

Suppose there exists  $\phi(x, y)$  s.t.

$$d\phi(x, y) = P(x, y)dx + Q(x, y)dy = 0, \quad (7.8)$$

and then the solution is  $\phi(x, y) = \text{constant}$ . Notice that since

$$P = \frac{\partial \phi}{\partial x}, Q = \frac{\partial \phi}{\partial y}, \quad (7.9)$$

it must be that since mixed partials commute,

$$\frac{\partial P}{\partial y} = \frac{\partial Q}{\partial x} \quad (7.10)$$

is a necessary and sufficient condition for the equation to be exact. In the language of differential forms, we may say that the equation is closed,

$$d\omega = Pdx + Qdy = 0, \quad (7.11)$$

but that closed does not imply exact,  $\omega = d\mu$  for some  $\mu$ . (Exact implies closed since  $d^2 = 0$ .)

<sup>11</sup>Strictly, this manipulation is justified by differential forms.

However, if the equation is not exact, we could think of multiplying by some unknown function  $\alpha(x, y)$  such that it becomes exact:

$$\alpha(x, y)P(x, y)dx + \alpha(x, y)Q(x, y)dy = 0. \quad (7.12)$$

In general there's no formula for finding such an  $\alpha$ , but there is one case where we can do this, namely the linear case. Let

$$y' + p(x)y = Q(x). \quad (7.13)$$

We then write

$$D(\alpha y) = \alpha' y + \alpha y', \quad (7.14)$$

If we multiply our linear equation by

$$\alpha(x) = \exp\left(+\int^x dx' p(x')\right) \quad (7.15)$$

then we claim the equation becomes exact (a total derivative), since

$$D(\alpha y) = p(x)\alpha y + \alpha y'. \quad (7.16)$$

Hence

$$D[\alpha y] = \alpha(x)Q(x), \quad (7.17)$$

so we can just integrate:

$$y(x) = \frac{1}{\alpha(x)} \left[ \int^x dx' \alpha(x')Q(x') + C \right]. \quad (7.18)$$

We can think of this as a the sum of a homogeneous solution and a particular solution, i.e. when  $Q(x) = 0$  we just get  $y(x) = \frac{C}{\alpha(x)}$  and there is a particular solution from dealing with  $Q(x) \neq 0$ .

**Second-order equations** Let us consider the case of second-order equations with constant coefficients,

$$a \frac{d^2 y}{dx^2} + b \frac{dy}{dx} + cy = f(x). \quad (7.19)$$

There are various ways of solving such equations. As before, we can solve the homogeneous equation and then add back in the particular solution turning on the source term. There are some methods in the literature such as the method of undetermined coefficients or variation of parameters.

We can denote our equation of order  $n$  by

$$\frac{d^n y}{dx^n} + a_{n-1} \frac{d^{n-1} y}{dx^{n-1}} + \dots + a_0 y = f(x), \quad (7.20)$$

or in the simplified notation

$$p_n(D)y = f(x), \quad p_n(D) = D^n + a_{n-1}D^{n-1} + \dots \quad (7.21)$$

Then let us note that derivatives obey the following identities<sup>12</sup>

$$\begin{aligned} D + \lambda &= e^{-\lambda x} D e^{\lambda x} \\ D - \lambda &= e^{\lambda x} D e^{-\lambda x} \\ (D - \lambda)^n &= e^{\lambda x} D^n e^{-\lambda x}. \end{aligned}$$

This is equivalent to writing that

$$e^{-\lambda x} D(e^{\lambda x} y) = y' - \lambda y. \quad (7.22)$$

Now the homogeneous second order equation takes the form

$$(aD^2 + bD + c)y = 0. \quad (7.23)$$

We know how to factor expressions. If we can factor the differential operator, then the solutions are straightforward because we know what  $D \pm \lambda$  does.

The nature of solutions (how nicely will the operator factor) splits into three cases depending on the discriminant  $\Delta = b^2 - 4ac$ . if  $\Delta > 0$  then the equation has two real roots  $\lambda_1, \lambda_2$ , so that

$$(D - \lambda_1)(D - \lambda_2)y = 0, \quad (7.24)$$

<sup>12</sup>That is, we are using the generator of translations and exponentiating.

with solution

$$y = Ae^{\lambda_1 x} + Be^{\lambda_2 x}. \quad (7.25)$$

If  $\Delta < 0$  then there are two complex roots  $\lambda_{\pm} = \mu + i\omega$ , so that the equation factors as

$$\left[(D - \mu)^2 + \omega^2\right]y = 0. \quad (7.26)$$

Recall that  $(D^2 + \omega^2)y = 0$  has solutions  $y = Ae^{i\omega x} + Be^{-i\omega x}$ , so then

$$\begin{aligned} p(D)y &= (e^{\mu x} D^2 e^{-\mu x} + \omega^2)y \\ &= e^{\mu x} (D^2 + \omega^2) e^{-\mu x} y \\ &= 0 \implies e^{-\mu x} y = Ae^{i\omega x} + Be^{-i\omega x}. \end{aligned}$$

Hence the solutions are both oscillating and growing or decaying exponentially with time,

$$y = e^{\mu x} (Ae^{i\omega x} + Be^{-i\omega x}) \quad (7.27)$$

Finally, we could have  $\Delta = 0$ , which gives a double root at  $\lambda$ . Then

$$\begin{aligned} p(D)y &= (D - \lambda)^2 y = 0 \\ &= e^{\lambda x} D^2 e^{-\lambda x} y = 0. \end{aligned}$$

But this tells us we want some  $z$  such that  $D^2 z = 0$ , and hence

$$e^{-\lambda x} y = A + Bx, \quad (7.28)$$

which gives

$$y = e^{+\lambda x} (A + Bx). \quad (7.29)$$

This is equivalent to the statement that  $e^{\lambda x}, xe^{\lambda x}$  are in the kernel of  $(D - \lambda)^2$ . In each of these cases, we look for the kernel of some differential operator and set it equal to a known expression.

Now can we do this for an inhomogenous equation? Indeed we can.

**Example 7.30.** Consider

$$y'' - 3y' + 2y = e^x, \quad (7.31)$$

so

$$(D^2 - 3D + 2)y = (D - 1)(D - 2)ye^x, \quad (7.32)$$

so now  $(D - 1)z(x)e^x$  has solution  $z(x) = Ae^x + e^x$  and so

$$(D - 2)y = Ae^x + xe^x. \quad (7.33)$$

Then

$$y = e^x - xe^x - Ae^x + Be^{2x}. \quad (7.34)$$

**Example 7.35.** Suppose we want to find a particular solution to

$$3y'' - 2y' + 6y = 5e^{3x}. \quad (7.36)$$

Then my ansatz might be that the solution should be in the form of some constant times  $e^{3x}$ . But alternately, we could invert the differential operator. We know how to do this. With

$$(3D^2 - 2D + 6)y = 5e^{3x}, \quad (7.37)$$

we can write

$$y = \frac{1}{3D^2 - 2D + 6}(5e^{3x}) = \frac{1}{3(3)^2 - 2(3) + 6}5e^{3x} = \frac{5}{27}e^{3x}. \quad (7.38)$$

In general for

$$p_n(D)y = Ae^{\lambda x}, \quad (7.39)$$

if  $p_n(\lambda) \neq 0$  then

$$y = \frac{1}{p_n(\lambda)} Ae^{\lambda x} \quad (7.40)$$

and if  $p_n(\lambda) = 0$  then

$$y = \frac{1}{p'_n(\lambda)} x A e^{\lambda x}. \quad (7.41)$$

This is known as the exponential input theorem.

**General linear 2nd order equation** Let's consider the general case,

$$y'' + P(x)y' + Q(x)y = 0. \quad (7.42)$$

This is generally hard to solve. It's a little easier if  $P(x), Q(x)$  are analytic except at a finite number of poles. If we wanted to solve the equation near a nonsingular (ordinary) point, we could expand  $P$  and  $Q$  as power series and solve locally.

Alternately, we might have a *regular singular point* where  $P(x)$  has no worse than a single pole ( $P(x) \sim 1/x$ ) and  $Q(x)$  has no worse than a double pole ( $Q(x) \sim 1/x^2$ ). Beyond this there are also *irregular singular points* or sometimes essential singularities.

**Example 7.43.** Consider the following differential equation:

$$x(x-1)y'' + [(1+a+b)x+c]y' + aby = 0, \quad (7.44)$$

the hypergeometric equation. This equation has regular singular points at  $x = 0, x = 1$ . But we could also analyze the point at  $x \rightarrow \infty$ , where if we replace  $x = \frac{1}{z}$  and write  $w(z) = y(1/z)$ , then with a bit of chain rule, our standard second order equation becomes

$$w'' + \left[ \frac{2z - P(1/z)}{z^2} \right] w' + Q(1/z)w = 0. \quad (7.45)$$

Hence the equation we were given has a regular singular point at  $x = \infty$ .

There's one more nice fact about general linear second-order equations, which is that we can generally get rid of the order  $y$  term. Set  $y(x) = \alpha(x)z(x)$  for some  $\alpha$ . In particular if we choose

$$\alpha(x) = \exp\left(-\frac{1}{2} \int^x dx' P(x')\right) \quad (7.46)$$

then the equation (exercise) becomes

$$z''(x) + I(x)z(x) = 0 \quad (7.47)$$

for some  $I$  given by

$$I(x) = Q(x) - \frac{1}{2}P'(x) - \frac{1}{4}P(x)^2. \quad (7.48)$$

This expression  $I$  is actually an invariant— if two different-looking differential equations have the same value of  $I$ , they are really the same equation.