1 Summary

This is an informal overview of the "system och signaler" course.

2 Linear algebra precursor

To fully motivate the definition of the fourier transform, we'll need a bit of linear algebra first.

2.1 Basis

I'll begin with the concept of a *basis*. A list of vectors $v_1, \ldots, v_n$ can be thought of as a basis if every other vector can be represented by the basis by uniquely scaling and adding them.

(a) This **is** a basis for $\mathbb{R}^2$ (b) This **is not** a basis for $\mathbb{R}^2$
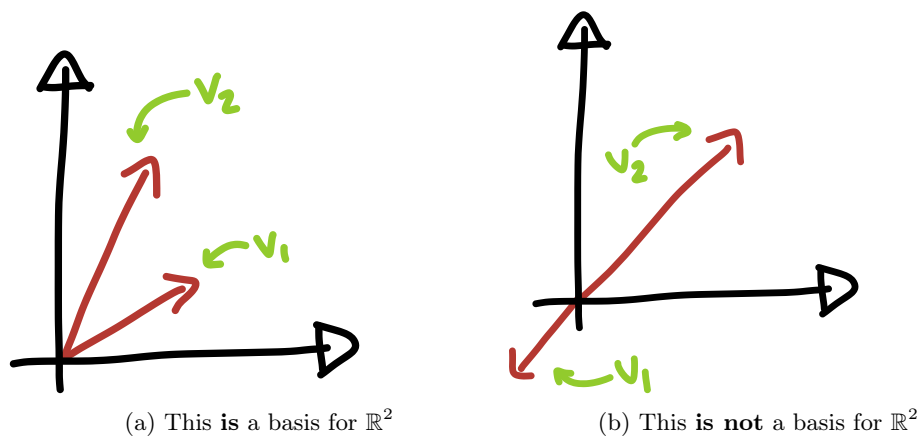
2.1.1 Example

For example the vectors

$$v_1 = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, v_2 = \begin{bmatrix} 0 \\ 1 \end{bmatrix}$$

is a basis for $\mathbb{R}^2$.

2.1.2 Question

Are the vectors

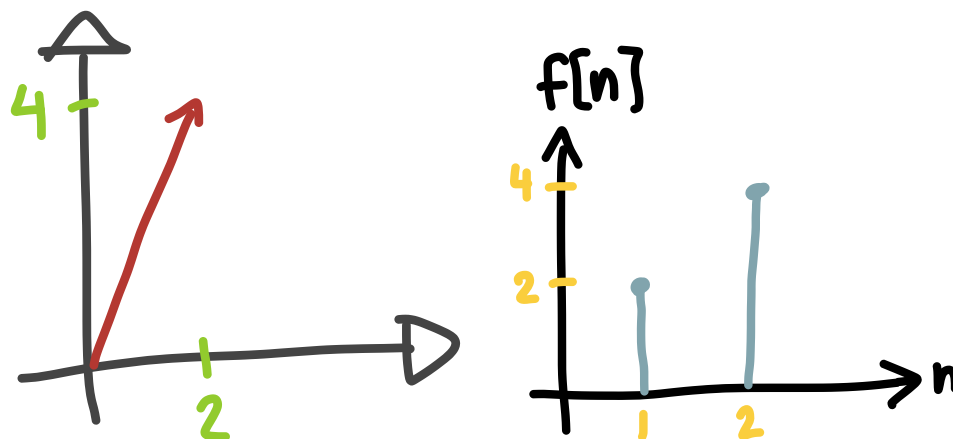$$v_1 = \begin{bmatrix} 1 \\ 0 \end{bmatrix}, v_2 = \begin{bmatrix} 2 \\ 0 \end{bmatrix}$$

a basis for $\mathbb{R}^2$? Why or why not?

2.2 Functions are vectors

From linear algebra we're used to thinking of vectors like

$$v = \begin{bmatrix} 2 \\ 4 \end{bmatrix}$$

as arrows, perhaps on the cartesian plane in the case of 2D vectors.


But an important insight we'll need on our journey into the fourier transform is that functions can actually be seen as vectors! To motivate this, let's just think about how we could represent the v above as a function f. Well the first entry has value 2, and the second entry has value 4, so

$$f[1] = 2, f[2] = 4$$

That certainly makes a lot of sense! And you could easily imagine how to convert arbitrary-dimensional vectors into functions.

Using this notation we can efficiently plot vectors which would otherwise be impossible, like 6-dimensional vectors!
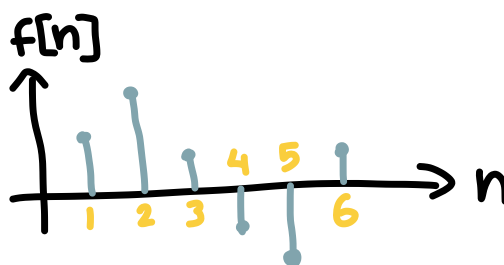
The whole point of this has been to illustrate that functions are indeed also vectors, which means it makes sense to define the dot-product between two functions, the length (aka norm) of a function, etc. In the next section we'll see how the dot-product of two functions could be defined.

(So far we've been working with discrete functions, but using similar arguments it's also possible to see that continuous functions are vectors.)

2.3 Dot-product ("skalärprodukt")

From linear algebra we learned that a dot-product between two real vectors (i.e. with all elements real valued)

$$v = \begin{bmatrix} a_1 \\ \vdots \\ a_n \end{bmatrix}, w = \begin{bmatrix} b_1 \\ \vdots \\ b_n \end{bmatrix}$$

is defined as

$$v \cdot w = a_1 b_1 + \cdots + a_n b_n$$

Note that $v \cdot v = a_1^2 + \cdots + a_n^2 = \|v\|^2$, since $\|v\|$ is defined as $\|v\| = \sqrt{a_1^2 + \cdots + a_n^2}$.

Above we restricted ourselves to vectors with real-valued entries, but what about vectors with complex-valued entries? What happends if we use the above definition without modification? Well suppose we want to find out the length of $u = \begin{bmatrix} i \\ 0 \end{bmatrix}$. Intuitively you'd think the length of this would be 1, i.e. $\|u\| = 1$. But look what happends if we try to calculate u using the fact that $\|u\| = \sqrt{u \cdot u}$

$$\|u\| = \sqrt{u \cdot u} = \sqrt{i^2 + 0^2} = i$$

So we get a complex length! Now that's not very intuitive!

Instead of using this definition, we'll modify it slightly to ensure the norm is always a positive real. Recall that if z is a complex number, then $z\bar{z}$ is always a positive real. Using this as inspiration we define the dot-product of two complex vectors (i.e. with complex-valued entries)

$$v' = \begin{bmatrix} z_1 \\ \vdots \\ z_n \end{bmatrix}, w' = \begin{bmatrix} h_1 \\ \vdots \\ h_n \end{bmatrix}$$

to be

$$v' \cdot w' = z_1 \overline{h_1} + \cdots + z_n \overline{h_n}$$

Notice that when v' and w' are real, this turns into the normal (real) dot-product (since $a = \bar{a}$, for real a). So we've really just extended the notion of the dot-product, not replaced it.

2.3.1 Dot-products of functions

So far we've only looked at the dot-product of vectors in $\mathbb{R}^n$ and $\mathbb{C}^n$, but it's actually possible to also define the dot-product of functions.

To motivate the definition, let's recall that for two complex vectors

$$v = \begin{bmatrix} v_1 \\ \vdots \\ v_n \end{bmatrix}, w = \begin{bmatrix} w_1 \\ \vdots \\ w_n \end{bmatrix}$$

the following is their dot-product

$$v \cdot w = v_1 \overline{w_1} + \cdots + v_n \overline{w_n}$$

note that this can also be written as

$$v \cdot w = \sum_{k=1}^{n} v_k \overline{w_k}$$

So the natural definition for a couple of discrete functions $t[x], h[x]$, where x goes from 1 to n, is just

$$t[x] \cdot h[x] = \sum_{k=1}^{n} t[k] \overline{h[k]}$$

What about continuous functions? We'll begin with periodic functions. When we go from the discrete to the continuous, series turn into integrals, thus it's be pretty natural that the definition should be something like

$$p(x) \cdot q(x) = \int f(x) \overline{g(x)} dx$$

where $p(x)$ and $q(x)$ are periodic with period T_0. Notice that I didn't write the integration limits. What should these be? Well since an integral during a period of a periodic function really contains all the information about the area we need, it'd be reasonable to have the limits be one of the periods (e.g. from 0 to T_0)

$$p(x) \cdot q(x) = \int_{T_0} f(x) \overline{g(x)} dx$$

Note that the integral will be equivalent no matter which period we choose.

What about aperiodic functions $f(x)$ and $g(x)$? These can be seen as periodic functions but with an infinite period, so we just extend the above to

$$f(x) \cdot g(x) = \int_{-\infty}^{\infty} f(x) \overline{g(x)} dx$$

Sometimes the dot-product between two functions is denoted $\langle f, g \rangle$, to avoid confusion from normal multiplication. I'll use this notation in the remainder of this document.

2.4 Orthagonality

Orthagonal is just a fancy word for perpendicular.

So the definition of two vectors v and w being orthagonal is just that

$$v \cdot w = 0$$

or, using the notation for the dot-product between function two functions f and g,

$$\langle f, g \rangle = 0$$

We're especially interested in orthogonal bases, i.e. where all vectors in the basis are pairwise orthogonal. Why do we care about these types of bases? To motivate this, let's investigate what happends when we have a basis which **isn't** pairwise orthogonal, like

$$v_1 = \begin{bmatrix} 1 \\ 2 \end{bmatrix}, v_2 = \begin{bmatrix} 2 \\ 1 \end{bmatrix}$$

4

Since $v_1 \cdot v_2 = 4 \neq 0$, this is not an orthagonal basis.

Now suppose we're given some vector

$$v = \begin{bmatrix} 4 \\ 0 \end{bmatrix}$$

and we want to find the scalars α_1, α_n such that

$$v = \alpha_1 v_1 + \alpha_2 v_2$$

In other words, such that

$$\begin{bmatrix} 4 \\ 0 \end{bmatrix} = \alpha_1 \begin{bmatrix} 1 \\ 2 \end{bmatrix} + \alpha_2 \begin{bmatrix} 2 \\ 1 \end{bmatrix} = \begin{bmatrix} \alpha_1 \\ 2\alpha_1 \end{bmatrix} + \begin{bmatrix} 2\alpha_2 \\ \alpha_2 \end{bmatrix}$$

So we need to solve the following system of equations

$$\begin{cases} \alpha_1 + 2\alpha_2 = 4 \\ 2\alpha_1 + \alpha_2 = 0 \end{cases}$$

How do we do this? You guessed it! Matrices!

$$\begin{bmatrix} 1 & 2 & | & 4 \\ 2 & 1 & | & 0 \end{bmatrix}$$

I'll skip the details, but eventually we get that

$$\alpha_1 = \alpha_2 = \frac{4}{3}$$

Which you should verify is correct.

Why did we do all this? To illustrate the point that arbitrary bases are annoying. What we really want, it turns out, are orthagonal bases!

2.4.1   Orthogonal bases

When we have an orthogonal basis it turns out there exists a really nice formula for each of the scalars. To motivate this, recall that the formula for projecting a vector v onto another vector u is denoted by $proj_u(v)$ and is given by

$$proj_u(v) = \frac{v \cdot u}{\|u\|^2} u$$

If u is normalized (i.e. $\|u\| = 1$) this turns just into

$$proj_u(v) = (v \cdot u)u$$

So with this in mind, let's return to the original problem. Suppose we have some vector v and we want to find the scalars $\alpha_1, \ldots, \alpha_n$ such that

$$v = \alpha_1 e_1 + \cdots + \alpha_n e_n$$

where $e_1, \ldots, e_n$ is an orthonormal (i.e. pairwise orthogonal and normalized) basis. Then we can sort of imagine "projecting" v onto all basis vectors. Thus the scalars should be

$$v = (v \cdot e_1)e_1 + \cdots + (v \cdot e_n)e_n$$

and indeed this is the correct representation of v in terms of the basis!

2.5 Can we use the projection formula for non-orthagonal bases?

Why can't we use the projection formula for non-orthagonal bases? Well, because it wouldn't give us the correct scalars. See the figure below for an illustration of where how fails.
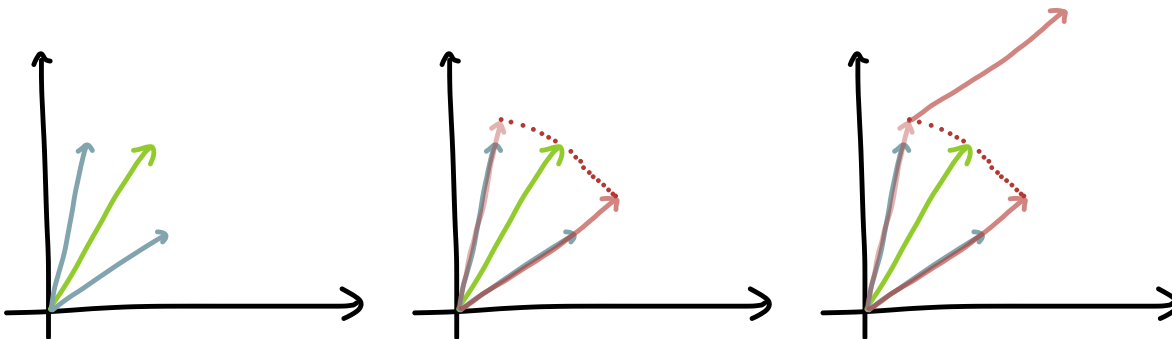
When adding the projections of the red vector onto the blue basis we just don't get back the red vector. Instead we would have to solve a system of equations. The whole effort could be avoided by using an orthogonal basis.

3 Fourier series

Finally we come to what the course is actually about. Luckily this step will be super simple thanks to the background we've provided.

3.1 Searching for an orthonormal basis to the periodic functions

So we've motivated why orthonormal bases are important/nice. And we've seen some examples of orthonormal bases for $\mathbb{R}^2$. We'd now like to find an orthonormal basis for the set of periodic functions.

3.1.1 A first attempt (sinusoids of the same phase)

So suppose we have some periodic function $x_{T_0}(t)$ with period T_0. What might be an appropriate choice of building blocks to represent this? Well some of the most famously periodic functions are the sinusoids, so it seems reasonable that summing a bunch of these will eventually construct our function

$$x_{T_0}(t) \overset{?}{=} \sum_{n=0}^{\infty} \alpha_n sin(n\omega_0 t)$$

Here ω_0 denotes the angular frequency, which is given by $\omega_0 = \frac{2\pi}{T_0}$ (since $\omega_0 = 2\pi f_0$ and $T_0 = \frac{1}{f_0}$).

Notice here that all of these sinusoids have the same phase (they all start at zero and end at zero). Is there any problems with this? Are there periodic functions which cannot be represented as sinusoids, all of the same phase? It turns out yes.

Consider the simple sine, shifted up by 1

$$s(t) = 1 + sin(t)$$

Of all periodic function, surely you'd expect this to be representable by a series of sines! But consider what happends at $t = 0$. Here $s(0) = 1$ but our sinusoids all start at zero! Adding an infinite amount of zeroes will never yeild 1, so this function is not representable by our series.
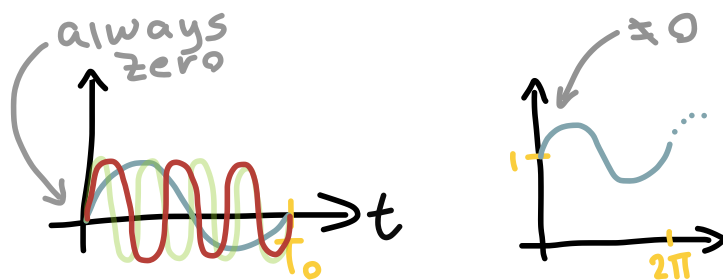
You might think; "well if we need the x(0) to be 1, perhaps that's a sign sines are a bad choice, switch to a series of cosines instead!" But the problem remains here, for a series of cosines (all of the same phase) always start at 1, so how do we represent the most basic function of all; $x(t) = 0$?

3.1.2 A second attempt (sinusoids of different phases)

So sinusoids of the same phase won't work. What about if we also allow the phase to vary?

$$x_{T_0}(t) \overset{?}{=} \sum_{n=0}^{\infty} \alpha_n sin(n\omega_0 t + \phi_n)$$

Making $\omega_0 = 1, \phi_0 = \pi, \alpha_0 = 1, \phi_1 = 0, \alpha_1 = 1$ and all other α's zero yields

$$s(t) = 1 \cdot sin(0 \cdot t + \pi) + 1 \cdot sin(1 \cdot t + 0) + 0 + \cdots = 1 + sin(t)$$

thus our problematic function above **will** be representable by this series. Seems promising.

   If you've worked with sinusoids before, however, you know it's annoying. There's trigonometric identities everywhere and integration is a pain. Although this series is promising, we'd really like to work with simpler terms. In the following section I'd like to argue that the series is actually equivalent to a much nicer series of complex exponentials.

3.1.3 A third attempt (compex exponentials, optional[1])

We begin by replacing all ϕ_k with $\theta_k + \frac{\pi}{2}$, thus transforming all sines into cosines. ϕ_k is just an arbitrary constant of our choosing so there's nothing stopping us from doing this.

$$\sum_{n=0}^{\infty} \alpha_n sin(n\omega_0 t + \frac{\pi}{2} + \theta_n) = \sum_{n=-\infty}^{\infty} \alpha_n cos(n\omega_0 t + \theta_n)$$

Now we invoke eulers formula

$$cos(n\omega_0 t + \theta_n) = \frac{e^{j(n\omega_0 t + \theta_n)} + e^{-j(n\omega_0 t + \theta_n)}}{2}$$

which, when replaced into the series, becomes

$$\sum_{n=0}^{\infty} \frac{\alpha_n}{2} \left(e^{jn\omega_0 t + j\theta_n} + e^{-jn\omega_0 t - j\theta_n} \right) = \sum_{n=0}^{\infty} \left(\frac{\alpha_n}{2} \left(e^{j\theta_n} e^{jn\omega_0 t} \right) + \frac{\alpha_n}{2} \left(e^{-j\theta_n} e^{-jn\omega_0 t} \right) \right)$$

We now split up the sum and replace $\frac{\alpha_k}{2} e^{j\theta_n}$ with β_k. (Notice that the complex conjugate $\overline{\beta_k} = \frac{\alpha_k}{2} e^{-j\theta_n}$.)
So we get

$$\sum_{n=0}^{\infty} \beta_n e^{jn\omega_0 t} + \sum_{n=0}^{\infty} \overline{\beta_n} e^{-jn\omega_n} = \sum_{n=0}^{\infty} \beta_n e^{jn\omega_0 t} + \sum_{n=-\infty}^{0} \overline{\beta_{-n}} e^{jn\omega_n}$$

Now we finally combine the sums again to get

$$\sum_{n=-\infty}^{\infty} \beta_n e^{jn\omega_0 t}$$

And we're done. (Note that for $n < 0$, $\beta_n = \overline{\beta_{-n}}$, and that $\beta_0 = \alpha_0$.)

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[1]Here we just show that we can represent the infinite sum of sinusoids as an infinite sum of complex exponentials. It can be skimmed if you don't care for long derivations.

3.1.4 Orthonormal

So where are we? We're trying to find an orthonormal basis for the set of periodic functions. We began with the hunch that sinusoids might be appropriate. Sinusoids are annoying but luckily we found that the these were equivalent (due to eulers formula) to a set of complex exponentials. We'd now like to provide some rationale as to why these are an orthonormal basis for the set of periodic functions, meaning we should be able to represent *any* periodic signal as a linear combination (a sum of scaled versions) of these complex exponentials. Not only that, but since they're an *orthonormal* basis, we'll have a really simple formula for caluclating what each scalar should be!

Unfortunately we won't be able to show that they're a basis[2], but we will show that they're orthonormal!

Before starting we'll need need to assume that the period is 1, thus $\omega_0 = 2\pi$ (this is just to simplify things, it's also possible to show for arbitrary periods[3]).

We will denote the n:th complex exponential by $e_n(t)$

$$e_n(t) = e^{j2\pi nt}$$

We begin by showing that they're orthagonal. (Recall that $\omega_0 = 2\pi$.)

Orthagonal We want to show that if $n \neq m$ then $\langle e_n, e_m \rangle = 0$. Suppose $n \neq m$

$$\langle e_n, e_m \rangle = \int_0^1 e^{j2\pi nt}\overline{e^{j2\pi mt}}dt = \int_0^1 e^{j2\pi(n-m)t}dt = \frac{1}{2\pi(n-m)}\left(e^{j2\pi(n-m)} - e^0\right) = 0$$

Normalized We begin with the fact that

$$\langle e_n, e_n \rangle = \int_0^1 e^{j2\pi nt}\overline{e^{j2\pi nt}}dt = \int_0^1 e^{j2\pi(n-n)t}dt = \int_0^1 1dt = 1$$

Recall that (by definition) $\|e_n\| = \sqrt{\langle e_n, e_n \rangle}$, thus $\|e_n\| = 1$, which is what we wanted.

3.2 The final result

Now that we know the complex exponentials are an orthonormal basis for the set of periodic signals, what does that mean? It means every periodic signal $x(t)$ can be represented as a linear combination of complex exponentials

$$x(t) = \sum_{n=-\infty}^{\infty} \alpha_n e^{j\omega_0 nt}$$

and since the basis is orthonormal, we know that we can write each scalar α_k as

$$\alpha_k = \left\langle x(t), e^{j\omega_0 kt} \right\rangle = \int_{T_0} x(t)\overline{e^{j\omega_0 kt}}dt = \int_{T_0} x(t)e^{-j\omega_0 kt}dt$$

And now we're done! We've successfully written a periodic signal in terms of a linear combination of complex exponentials. In the next section we'll extend this idea to signals which *aren't* periodic.

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[2]There exists a proof, I've just never seen it. . .

[3]I'm lying here. . . Kind of. The problem is that there are two choices of basis functions. The first choice is orthonormal and the other is just orthogonal. TSDT84 uses the orthagonal basis, and I'm trying to stay true to the course, but the formulas and motivation become much cleaner/easier with the orthonormal case. The result is me trying to combine the two with lies.

4 Fourier transform

The idea behind the fourier transform is really the same as fourier series, but instead of limiting ourselves to periodic signals, we'll also be able to deal with non-periodic signals (with one caveat, that their norm has to be finite, but we'll get to that shortly).

Let's get right into it!

We originally learned that since the complex exponentials is an orthonormal basis for the set of periodic functions (signals), that for every periodic signals $x(t)$ there exists a set of coefficients α_n, where n ranges from $-\infty$ to ∞, such that

$$x(t) = \sum_{n=-\infty}^{\infty} \alpha_n e^{j\omega_0 n t}$$

In particular (since it's an *orthonormal* basis), that each term α_n could be written as[4]

$$\alpha_n = \left\langle x(t), e^{j\omega_0 n t} \right\rangle = \int_{T_0} x(t) e^{-j\omega_0 n t} dt.$$

We now want to extend this notion to aperiodic functions. To do this, we simply[5] let the period $T_0 \to \infty$. Remember that $T_0 = \frac{1}{f_0} = \frac{2\pi}{\omega_0} \implies \omega_0 = \frac{2\pi}{T_0}$, so as $T_0 \to \infty, \omega_0 \to 0$. We're kind of getting more and more terms in the series representation of $x(t)$, until we're essentially dealing with a continuum of ω's, and the sum turns into an integral, called the *inverse fourier transform*,

$$x(t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} \alpha_\omega e^{j\omega t} d\omega.$$

(The $\frac{1}{2\pi}$ factor is just something we engineers like because it makes some identities easier. The "mathematicians" fourier transform does not have the factor[6].) where α_ω is the *fourier transform*, denoted $\alpha_\omega = \mathscr{F}\{x(t)\}(\omega)$, and is defined (as you'd expect)

$$\mathscr{F}\{x(t)\} = \left\langle x(t), e^{j\omega t} \right\rangle = \int_{-\infty}^{\infty} x(t) e^{-j\omega t} dt$$

(Recall that the negative exponent is due to taking the conjugate of the second factor, which we motivated earlier.)

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[4]I feel I must remind the reader that using the definition of the dot-product as defined here is not really valid as it will not make the complex exponentials we've chosen an ortho*normal* basis. The proper definition should include a $\frac{1}{T_0}$ factor in front, but motivating the reason why complicates the concept of the dot-product and would take away from the main point this (informal) section is trying to make.

[5]Formally, the integral doesn't really make sense when we let $T_0 \to \infty$, but remember that we're talking informally here.

[6]This is a consequence of the different choices of basis functions mentioned above.

4.1 Convergence and finite length

Unfortunately there's one caveat with the fourier transform, and that is that only functions with finite length can be transformed, because only these will make the integral converge. Recall that the length of a vector/function $f(t)$ is defined as

$$\|f(t)\| = \sqrt{\langle f(t), f(t)\rangle} = \sqrt{\int_{-\infty}^{\infty} |f(t)|^2 dt}$$

Note that in the book the requirement is that the squared length $\|f\|^2$ be finite. This is done simply because then we don't have to deal with the square root. But of course these requirements are equivalent. (If the length is infinite then the squared length is infinite, and vice versa.)

Why make this restriction? So what if the fourier transform diverges! Well suppose we had a two functions $g(t)$ and $h(t)$ with infinite energy and $g(t) \neq h(t)$, thus

$$\mathscr{F}\{g(t)\} = \mathscr{F}\{h(t)\} = \infty$$

But how do we go about inverting this? Should $\mathscr{F}^{-1}\{\infty\} = g(t)$ or should $\mathscr{F}^{-1}\{\infty\} = h(t)$? Clearly we cannot have it both ways. Instead we avoid the problem by just not letting the transform be defined for these types of functions.

In the next section we'll investigate an way of dealing with these types of functions.

5 Laplace transform

There are two types of laplace transforms. The unilateral and the bilateral. We'll begin with the bilateral.

5.1 The bilateral laplace transform

The bilateral laplace transform $\mathscr{L}_{II}$ of a signal $x(t)$ is

$$\mathscr{L}_{II}\{x(t)\} = \int_{-\infty}^{\infty} x(t)e^{-st}dt$$

where s is the complex, dependent variable of the laplace. It can be seen as a generalization of the fourier transform. To see the motivation behind this statement, we have to dig into what the definition really means.

Since s is complex, let's rewrite it into $s = \sigma + j\omega$. What do we get?

$$\int_{-\infty}^{\infty} x(t)e^{-st}dt = \int_{-\infty}^{\infty} x(t)e^{-(\sigma+j\omega)t}dt = \int_{-\infty}^{\infty} x(t)e^{-\sigma t}e^{-j\omega t}dt$$

But this looks suspiciously like the fourier transform of $x(t)e^{-\sigma t}$. Why would we ever want to do that? What's the point of all this?

Recall what the central problem with the fourier transform was. We cannot analyze signals using the fourier transform if they have infinite energy. A simple example of this is the signal (called the "unit step" function)

$$u(t) = \begin{cases} 1 \text{ if } t \geq 0 \\ 0 \text{ if } t < 0 \end{cases}$$

which has energy

$$\langle u(t), u(t) \rangle = \int_{-n}^{n} u(t)dt = \int_{0}^{n} 1dt \to \infty \text{ as } n \to \infty$$

thus $u(t)$ is not analyzable with the fourier transform. But can we analyze it with the laplace transform? Yes!
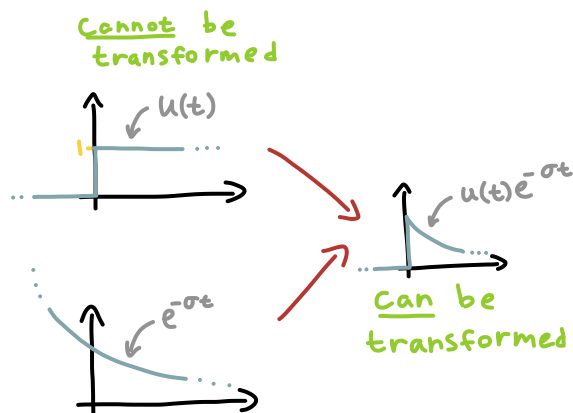
The power of the laplace is that uses an exponentially decreasing function $e^{-\sigma t}$ to sort of "push down" signals which have infinite energy into finite energy.

5.2 The unilateral laplace transform

The unilateral laplace transform, denoted $\mathscr{L}_I$ or just $\mathscr{L}$, of a signal $x(t)$ is

$$\mathscr{L}\{x(t)\} = \int_{0^-}^{\infty} x(t)e^{-st}dt$$

where 0^- means we also include the value of $x(t)$ at $t = 0$, which becomes important when you laplace transform the (or functions containing the) dirac.

Following analogy of the bilateral laplace, the unilateral laplace can be seen as taking the signal $x(t)$, then just ignoring the parts where $t < 0$.

5.3 Region of convergence

Some care should be taken when evaluating the laplace. The main problem is that it (usually) doesn't converge for all possible choices of s. Instead we get a "region" where s will make the transform converge to a useful value (i.e. not diverge into $\pm\infty$). This is called the "*region of convergence*", abbreviated ROC.

5.3.1 Example

Let's take an example to illustrate this. Consider what happends when we try to evaluate the bilateral[7] laplace transform of $u(t)$. We'll assume $s \neq 0$ and then consider $s = 0$ in a separate case below

$$\mathscr{L}_{II}\{u(t)\} = \int_{-\infty}^{\infty} u(t)e^{-st}dt = \int_{0}^{\infty} e^{-st}dt = \frac{e^{-st}}{-s}\Big|_{t=0}^{t=\infty} = \left[s = \sigma + j\omega\right] = -\frac{1}{s}\left(e^{-\sigma t}e^{-j\omega t}\Big|_{t=0}^{t=\infty}\right) \quad (1)$$

Now let's consider the upper and lower limits separately. We begin with the simple case; $t = 0$.

$$e^{-\sigma t}e^{-j\omega t}\Big|_{t=0} = e^0 e^0 = 1$$

What about $t = \infty$? The intuitive way[8] of thinking about this is; $e^{-\sigma t}e^{-j\omega t}$ is a complex number $(re^{j\phi})$. As $t \to \infty$, $e^{-j\omega t}$ just spins round and round, what really determines whether the limit diverges or converges is how the length, $e^{-\sigma t}$, behaves as $t \to \infty$, thus we can ignore the second term and instead just think about $e^{-\sigma t}$ as $t \to \infty$.

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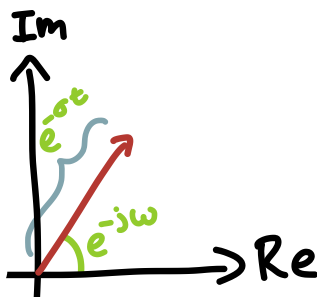
[7]Note that the choice of bilateral vs unilateral laplace transform in this case is arbitrary, the bilateral laplace quickly turns into the unilateral one. This is true in general when the signal is causal (i.e. when $x(t) = 0$ for $t < 0$)

[8]The rigorous way is not part of the course, and requires complex analysis I suppose.
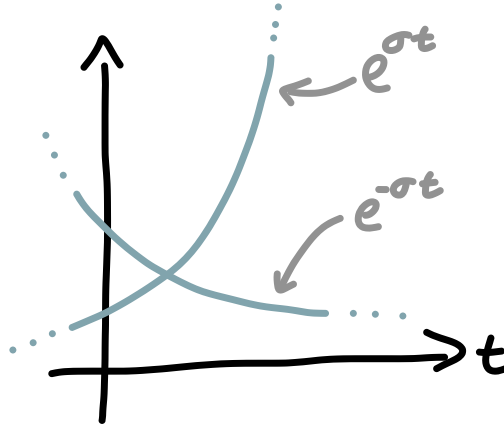
Now think about[9] what happends if $\sigma > 0$ and $\sigma < 0$. For the case of $\sigma > 0$, $e^{-\sigma t}$ becomes exponentially decreasing, which will converge to 0 as $t \to \infty$. What about $\sigma < 0$? Here it becomes an exponentially increasing, so very divergent!

Here's where the region of convergence comes in. We're interested in finding for what values of s the transformation converges. We just found that it will only converge for $Re\{s\} = \sigma > 0$, thus this will be our ROC!

Now let's continue with (1). We'd found that (still assuming $s \neq 0$)

$$\mathcal{L}_{II}\{u(t)\} = -\frac{1}{s}\left(e^{-\sigma t}e^{-j\omega t}\Big|_{t=0}^{t=\infty}\right) = -\frac{1}{s}(0-1) = \frac{1}{s}$$

Provided $\sigma > 0$.

What if $s = 0$?

$$\mathcal{L}_{II}\{u(t)\}(0) = \int_{-\infty}^{\infty} u(t)e^{-0 \cdot t}dt = \int_{0}^{\infty} dt \to \infty$$

Clearly $s = 0$ is not in the region of convergence of $u(t)$. This was already included in the $Re\{s\} > 0$ finding, so not terribly exciting.

5.3.2   Final result

The final result we got was that

$$\mathcal{L}_{II}\{u(t)\} = \frac{1}{s}$$

with a region of convergence of $Re\{s\} > 0$.

---

[9]I conveniently ignored the case of $\sigma = 0$. In this case we'll need to return to the $e^{-j\omega t}$ term, since $e^{-\sigma t}\Big|_{\sigma t} = e^0 = 1$. The motivation behind ignoring this is that $e^{-j\omega t}$ will just spin around and round and, just like $sin(t)$, will not converge as $t \to \infty$, thus $\sigma = 0$ is not included in the ROC.

5.4   Why is it called a region?

So we got that the region of convergence for $\mathscr{L}_{II}\{u(t)\}$ to be $Re\{s\} > 0$. But why do we call it a *region* of convergence? It seems more appropriate to call it an "interval" of convergence! But remember that s is a complex number, and that we're only restricting the real part of s. Think about the whole complex plane. How would we visualize the ROC?
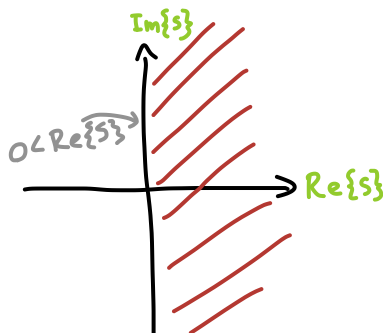
As you see in the picture above, although the ROC can really be described as an interval on the real line, in the complex plane it's really a region.

5.5   Relation between laplace and fourier transform

Earlier we said that the laplace can be seen as a generalization of the fourier transform. Now that we know what the ROC is, I'd like to briefly return to this concept. Recall that

$$\mathscr{L}_{II}\{x(t)\} = \int_{-\infty}^{\infty} x(t)e^{-st}dt = \left[s = \sigma + j\omega\right] = \int_{-\infty}^{\infty} \left(x(t)e^{-\sigma t}\right)e^{-j\omega t}dt = \mathscr{F}\{x(t)e^{-\sigma t}\}$$

If we now let $\sigma = 0$ we get

$$\mathscr{F}\{x(t)\} = \mathscr{L}_{II}\{x(t)\}\Big|_{Re\{s\}=0}$$

It's important to realize that this wonderful little relation is **only** true if $Re\{s\} = 0$ is in the region of convergence.

Visualizing it on the complex plane, we can see the fourier transform as the imaginary axis! We let $Re\{s\} = 0$ then plug in some value $\omega = Im\{s\}$ on the horizontal axis (the imaginary axis) to get the value of the fourier transform at that point!
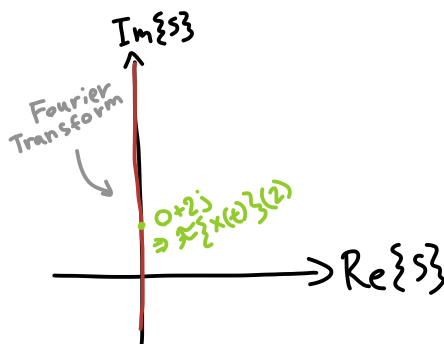
5.6   Poles and Zeros

Here are some laplace transforms that we might encounter

$$\frac{1}{s+1}, \frac{s+2}{(s-1)^2(s+3)}, \frac{s-5}{s^2(s+2)}, \cdots$$

As you can see, it's common that they can be represented as some polynomial divided by another polynomial.

The roots of the numerator are called the zeroes (i.e. where it is identically zero), and the roots of the denominator are called the poles, and are the points where the polynomials explode into $\pm\infty$[10]. These points are often drawn together with the ROC on the "s-plane" (which we've drawn above).

For instance $sin(t)u(t)$ has laplace transform $\frac{1}{s^2+1}$, with ROC $Re\{s\} > 0$. The polynomial has a pole at $s = \pm i$, thus the "s-plane" can be visualized as
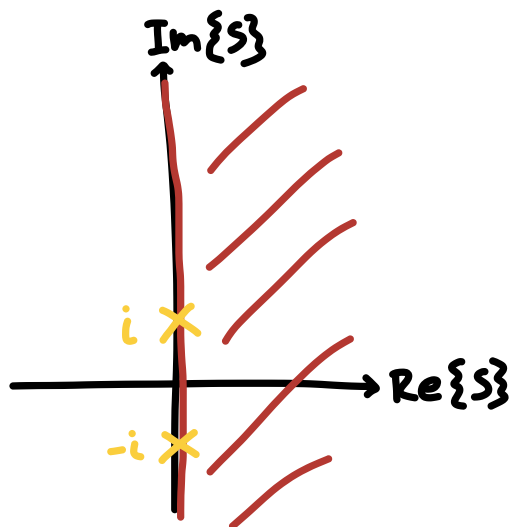
I won't prove it, but it turns out that the ROC will always have zeroes and/or poles at the edges.

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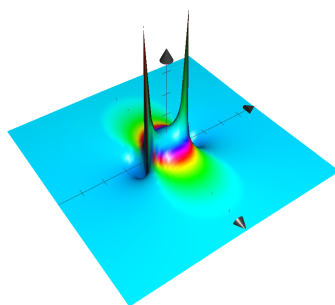
[10]The reason I say $\pm$ is because it will converge differently depending on which path you take on the complex plane. Though you could say that the poles are where the absolute value of the rational function explode to ∞.

5.7 Warnings

Here's a few common misconceptions and things to think about regarding the laplace.

5.7.1 Signals that are fourier transformable but not laplace transformable

Although we said the laplace could be seen as a generalization of the fourier transform, it's important to understand that there are some functions which are fourier transformable but **not** laplace transformable!

The canonical example is $\mathscr{L}_{II}\{1\}$

$$\mathscr{L}_{II}\{1\} = \int_{-\infty}^{\infty} e^{-st}dt = \left[s = \sigma + j\omega\right] = \int_{-\infty}^{\infty} e^{-\sigma t}e^{-j\omega t}dt$$

For $\sigma < 0$ the right side grows to infinity as $t \to \infty$, and for $\sigma > 0$ the left side grows to infinity as $t \to -\infty$.
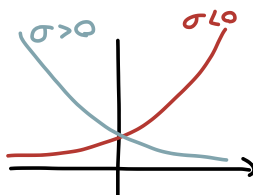

But we conveniently ignored the case of $\sigma = 0$. Well here it also won't converge. But remember this particular example wasn't a problem for the fourier transform. We just used diracs instead.

Essentially anything whose fourier transform contains a dirac, *isn't* laplace transformable (because "we just don't involve the dirac for the laplace" according to the lecturer).

5.7.2 Complex heightmap

The visualization of the laplace transform, displaying the ROC can be very useful. But it's important to realize that what we're viewing is not a heightmap, i.e. we can't make the plot 3D to visualize the "actual" transform, because each point is a complex number.

Let's take $\mathscr{L}\{u(t)\} = \frac{1}{s}$, with ROC $Re\{s\} > 0$ as an example. In order to fully see what value the transformation has at a given point, we need two plots, one for the real part and one for the imaginary part.
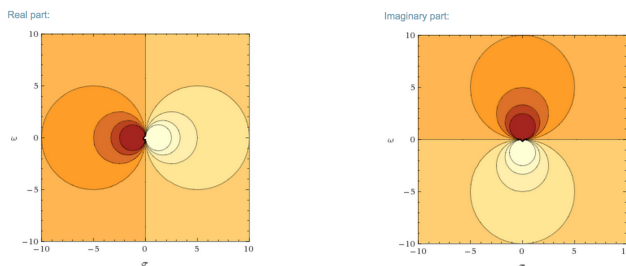

Note that we're displayin the whole plane here, whereas the transformation will only converge for $Re\{s\} > 0$. The image may show a finite value for $Re\{s\} < 0$, but that does not mean the the transformation would converge there. $\mathscr{L}\{u(t)\} = \frac{1}{s}$ is only valid for $\sigma = Re\{s\} > 0$.

6 Brief interlude

Where are we? We started investigating periodic, continuous, signals. We showed that as long as these have finite energy, then we could write these as a sum (linear combination) of complex exponentials. We then moved on to aperiodic signals and showed that by letting the period tend to infinity, these could also be analyzed with the *fourier transform*.

We took a quick sidestep investigating signals with infinite energy, and showed how these could be managed with the *laplace transform*.
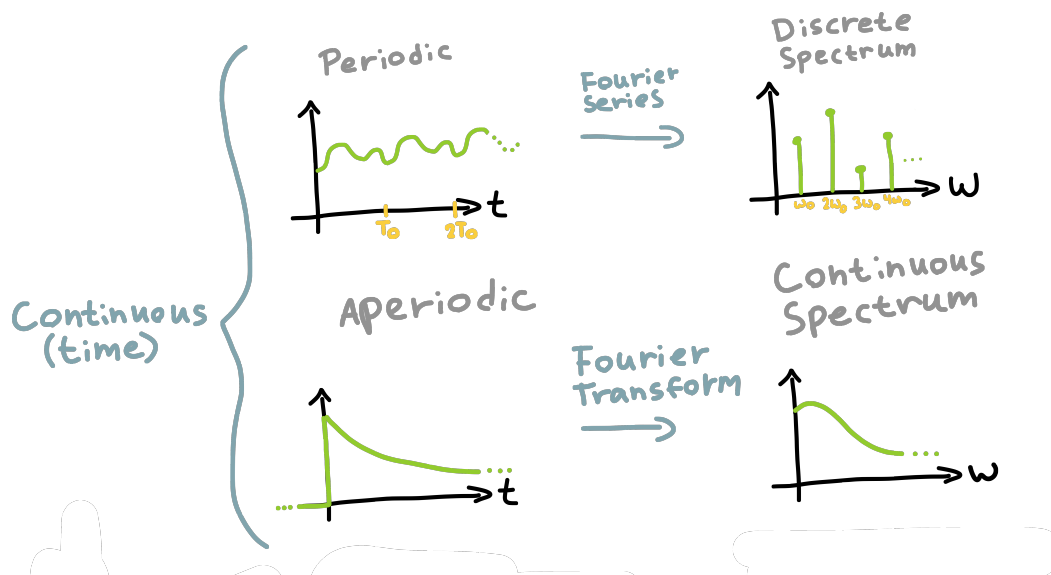
Now we're ready to leave the contiuous world and enter the discrete, starting with the z-transform and then moving on with the discrete-time fourier transform.

7 Sampling

Let's say we have a continous function $x(t)$ and we'd like to analyze it with a computer. Computers can only deal with discrete values, so what do we do? We have to find some way of discretizing the signal. This is called sampling.

The most obvious way of sampling the signal is just to periodically, say with period T, pluck out values and call this our sampled signal, let's call it $x[n] = x(nT)$, where n is an integer. Plug in $n = 3$ and we get the third sample-value of $x(t)$.

We'll get to how to choose the sampling-period T later when we get to the Nyquist–Shannon sampling theorem.

18

8 Z-transform

So far we've only been able to deal with (periodic and aperiodic) **contiuous** signals. Now we'd like to focus on the more practical, realistic, time-discrete signals. We call them "time-discrete" to emphasize the fact that it's the time-domain that's discrete, not the frequency-domain (like we had in the fourier series case).

8.1 Discrete signals in a continuous world

In the previous section we described how to sample a function; just create a new function $x[n] = x(nT)$. Unfortunately our tools are useless for dealing with this function, since we can only handle the contiuous. Luckily there's a way of constructing this sort of discrete function in the continous world. Just place a bunch of scaled deltas where the discrete values would be!
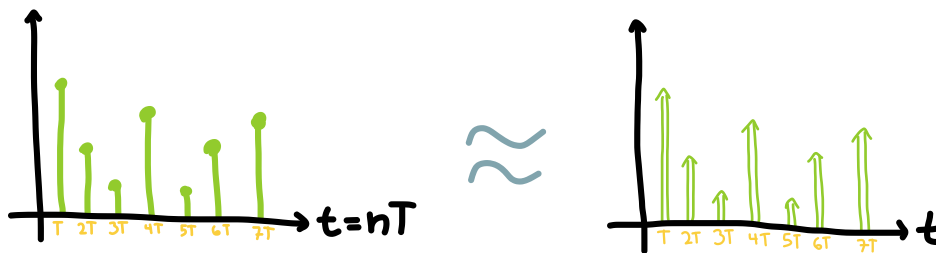
Sure, that sounds neat and all, but how would we go about representing this mathematically?

Well perhaps we can make use of the (conveniently named) *sampling property* of the dirac delta!
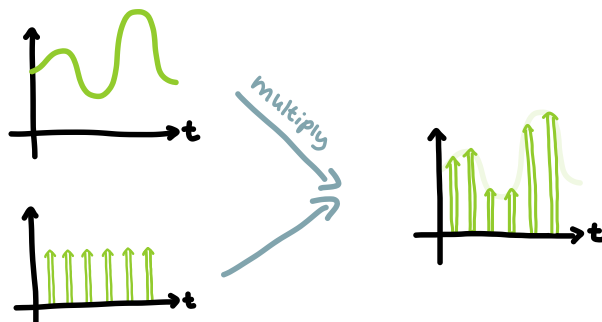
$$x(t)\delta(t - T) = x(T)\delta(t - T)$$

Now I'd like to introduce a new function, the III_T function (pronounced "Sha", a.k.a. the dirac comb or the dirac train). This is just an infinite amount of diracs spaced T apart.

$$\mathrm{III}_T(t) = \sum_{n=-\infty}^{\infty} \delta(t - nT)$$

Inspired by the sampling property, look what happends if we multiply III_T with $x(t)$

$$\mathrm{III}_T x(t) = \sum_{n=-\infty}^{\infty} x(t)\delta(t - nT) = \sum_{n=-\infty}^{\infty} x(nT)\delta(t - nT) = \sum_{n=-\infty}^{\infty} x[n]\delta(t - nT)$$

Let $\overline{x}(t) = \mathrm{III}_T x(t)$ and we've got our function!

8.2 Converting continuous transforms into discrete ones

So now that we know how to construct "discrete" signals in the contiuous world, it's easy to just convert our existing transforms into discrete equivalents. We'll begin with the laplace transform and we'll call its discrete equivalent the *z-transform*.

Let's just laplace transform $\overline{x}(t)$

$$\mathcal{L}_{II}\{\overline{x}(t)\} = \int_{-\infty}^{\infty} \overline{x}(n)e^{-st}dt$$

$$= \int_{-\infty}^{\infty} \left(\sum_{n=-\infty}^{\infty} x[n]\delta(t - nT) \right) e^{-st}dt$$

$$= \sum_{n=-\infty}^{\infty} x[n] \int_{-\infty}^{\infty} \delta(t - nT)e^{-st}dt$$

$$= \sum_{n=-\infty}^{\infty} x[n]e^{-snT} = \left[z = e^{sT} \right]$$

$$= \sum_{n=-\infty}^{\infty} x[n]z^{-n}$$

$$\overset{\text{def}}{=} \mathcal{Z}\{x[n]\}$$

Thus for discrete signals we'll use the z-transform, but remember that it's really the laplace in disguise.

8.3 Region of convergence

We've seen that the z-transform of a discrete signal is equivalent (by definition) to the laplace of a bunch of (equally spaced) diracs, thus in order to gain further insight into the z-transform we can just analyze the laplace.

There's one caveat though. Remember that the laplace transform depended on the variable s, whereas the z-transform depends on z. The relationship between them is $z = e^{sT}$. Thus when you take your signal, sample it, and wish to laplace transform it as if it was the original (continous) signal (using the z-transform), you must remember that everything will be in polar coordinates. Lines turn into circles.

Let's take an example to illustrate the point.

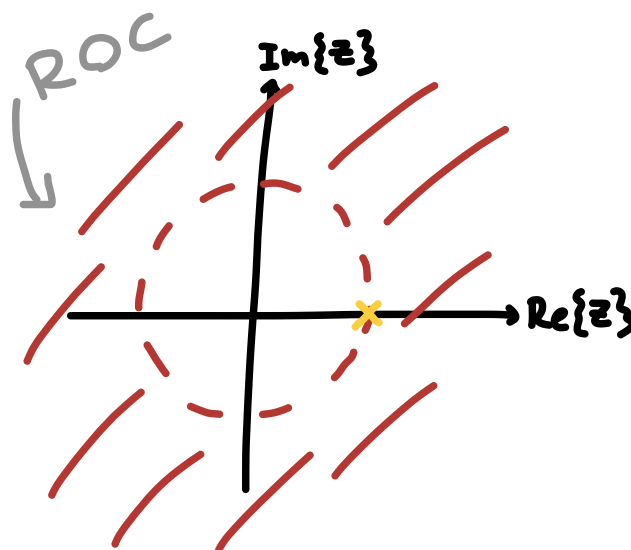### 8.3.1 Example

Say we want to find the z-transform of $u[t]$.

$$\mathscr{Z}\{u[n]\} = \sum_{-\infty}^{\infty} u[n]z^{-n} = \sum_{0}^{\infty} z^{-n}$$

Here's where the ROC comes in. We already know (from single variable calculus) that this converges to

$$\frac{z}{z-1}$$

if[11] $|z| > 1$. Also note that we have a pole at $z = 1$.

Put together this transform can be visualized in the "z-plane" as

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[11]and only if?

9   Discrete-time fourier transform

There's two ways of deriving the discrete-time fourier transform (DTFT).

- Convert discrete signals into contiuous equivalents (by multiplying with the dirac comb, thus sampling) and then apply the normal (continous) fourier transform which will (hopefully) produce an expression dependent on a discrete signal $x[n]$, which we'll take as the definition. This is how we did it when we went from the laplace to the z-transform.

- Alternatively, we could recognize the relation between the z-transform and the laplace, and the relation between the laplace and fourier, and try to find the definition from there instead.
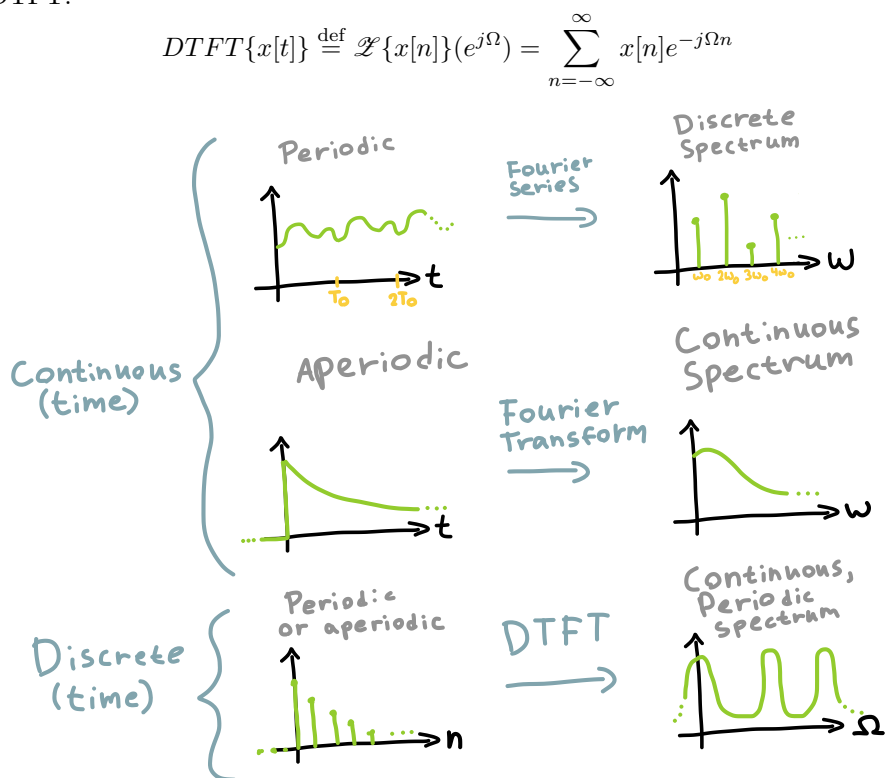
We'll be taking the second route.

10   Deriving the DTFT

So suppose we have some signal $x(t)$. We begin by sampling it $\overline{x}(t) = \text{III}_T x(t)$. Now recall that

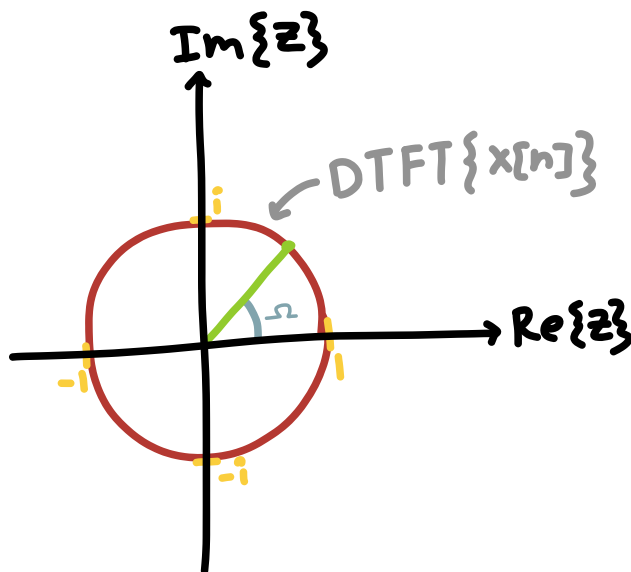$$\mathscr{F}\{\overline{x}(t)\} = \mathscr{L}\{\overline{x}(t)\}\Big|_{\sigma=0} = \mathscr{Z}\{x[t]\}(e^{-j\omega T})$$

assuming $Re\{s\} = 0$ is in the ROC. We now just rename ωT to Ω (called the "normalized" frequency) and we've got our DTFT!

$$DTFT\{x[t]\} \stackrel{\text{def}}{=} \mathscr{Z}\{x[n]\}(e^{j\Omega}) = \sum_{n=-\infty}^{\infty} x[n]e^{-j\Omega n}$$

Notice the duality between fourier series and the DTFT.

10.1   Why is the frequency spectrum of the DTFT periodic?

The standard way of explaining why the DTFT is periodic starts with a picture.

   As is clear from the derivation above, the values of the DTFT exist on the unit circle in the "z-plane" (assuming it's in the ROC, of course). This is analogous to how the fourier transform can be seen as the imaginary axis in the "s-plane". As you vary the angle Ω around the circle you'll eventually come back to where you started, which is why it's periodic.

10.1.1   Alternative (optional)

That's the usual way of explaining it, but I'd like to show it from a slightly different direction. Instead of taking the z-transform route we'll go directly from the fourier transform of the sampled signal $\overline{x}(t) = \text{III}x(t)$. We assume sampling period[12] $T = 1$ for simplicity, but the result is analogous for arbitrary periods.

$$DTFT\{x(t)\} = \mathscr{F}\{\overline{x}(t)\} = \mathscr{F}\{\text{III}x(t)\}$$

Now use the convolution/multiplication duality of the fourier transform and the fact[13] that $\mathscr{F}\{\text{III}\} = \text{III}$

$$\mathscr{F}\{\text{III}x(t)\} = \mathscr{F}\{\text{III}\} * \mathscr{F}\{x(t)\} = \text{III} * \mathscr{F}\{x(t)\}$$

Since III convolved with any function will make it periodic, we're done! What we've shown is that the DTFT spectrum is precisely[14] a version of the original signal, but periodized!
   Fun fact; this is one of the steps in the Nyquist-Shannon theorem proof! Just "unperiodize" it with a rect function to get

$$\mathscr{F}\{x(t)\} = \Pi(\text{III} * \mathscr{F}\{x(t)\})$$

Taking the inverse fourier transform of both sides yeilds the Nyquist-Shannon formula!

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[12]When we write III we really mean III_1.

[13]Provable using the Poisson summation formula.

[14]For period T the result would be $DTFT\{x(t)\} = \frac{1}{T}\text{III}_{\frac{1}{T}}\mathscr{F}\{x(t)\}$.

11 Systems

We now shift our focus to systems. A system is just a fancy word for "*function*". It takes as input functions and spits out new functions. An example of a system is

$$\mathcal{H}\{x(t)\} = 2 \cdot x(t)$$

If we're in a more mathematical context, sometimes we say *operator* instead of system. We'll begin by studying "continuous" systems (i.e. whose input are continuous functions). These could be implemented by an electric circuit. But later we'll also get to "discrete" systems (i.e. whose input is functions defined on discrete points), which could be implemented using digital circuits or in a computer program.
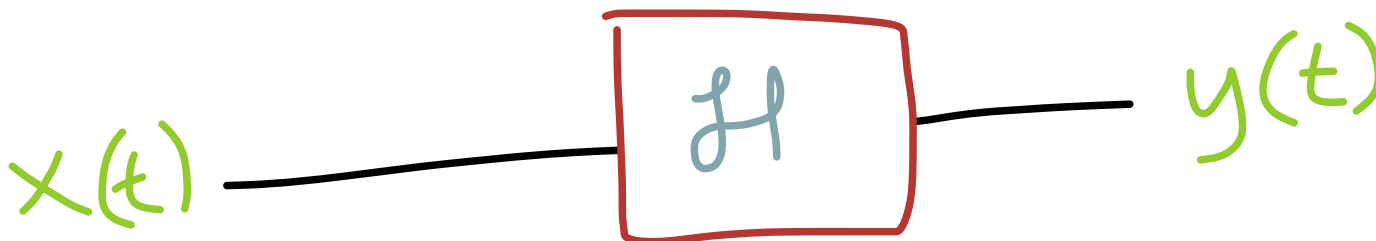
There are various properties a system can have. Among the ones we'll be primarily interested in are

- Time-invariant.
- Linear.
- Causal.
- Stable.
- Energy-free.

We'll get into these in depth now.

12 Time-invariance

The assumption on a system to be time-invariant just means that if we plug in $x(t)$ into our system $\mathcal{H}$ today, and get output $\mathcal{H}\{x(t)\} = y(t)$, then we're guaranteed to get the same output if we plug in the same input tomorrow $\mathcal{H}\{x(t+T)\} = y(t+T)$.

This is basically the same as requiring the system to be *state-less*. Of course it's not true for all systems we may want to investigate (e.g. circuits with state).

13 Linearity

That a system $\mathscr{H}$ is linear means specifically the following

- $\mathscr{H}$ is **homogenious**. In other words, $\mathscr{H}\{\alpha x\} = \alpha\mathscr{H}\{x\}$, for all $\alpha \in \mathbb{C}$, and all inputs x. (if $\mathscr{H}$ was defined on $\mathbb{R}^k$, this would have been all $\alpha \in \mathbb{R}$).

- $\mathscr{H}$ is **additive**. In other words, $\mathscr{H}\{x+y\} = \mathscr{H}\{x\} + \mathscr{H}\{y\}$, for all inputs x and y.

The reason we care about this property is because we get to use a lot of the theorems of linear algebra to understand our systems, as we'll see in the next section.

13.1 Linear operators are totally determined by its effect on a basis

The reason linearity is powerful is because we can totally understand a linear system by its effect on a basis.

Suppose we hade some linear operator $L : \mathbb{C}^k \to \mathbb{C}^k$, and $b_1, \ldots, b_k$ is some basis of $\mathbb{C}^k$. Now suppose we operate L on each basis vector, and get the following as a result

$$Lb_1 = v_1, \ldots, Lb_k = v_k$$

I claim that this information totally determines the effect of L on any vector $v \in \mathbb{C}^k$.

The proof is simple. Suppose $v \in \mathbb{C}^k$, is arbitrary. Since $b_1, \ldots, b_k$ is a basis, we can write v as a (unique) linear combination of this basis.

$$v = \alpha_1 b_1 + \cdots + \alpha_k b_k$$

Now let's plug in v into L again

$$Lv = L\{\alpha_1 b_1 + \cdots + \alpha_k b_k\} = \alpha_1 L\{b_1\} + \cdots + \alpha_k L\{b_k\} = \alpha_1 v_1 + \cdots + \alpha_k v_k$$

Since we know $\alpha_1, \ldots, \alpha_k$ and $v_1, \ldots, v_k$, we know the output of the system. And thus the proof is complete.

Notice that this assumed a finite dimensional vector space. Rest assured, however that the proof for this in the infinite-dimensional case is also possible, though it would require some extra assumptions (I think assuming absolutely integrable is enough).

14 LTI (linear and time-invariant)

So far we've only discussed what happends if we require our systems to be either linear or time-invariant, but some very interesting consequences come out of requiring our system to be both linear and time-invariant at the same time. This is often called an "LTI" system.

14.1 An LTI system is totally determined by its effect on the dirac

Say we had an LTI system $\mathscr{H}$. Since we know $\mathscr{H}$ is linear, in order to fully understand the system analytically, given arbitrary inputs, we may want to observe its effect on a basis. Suppose the basis we picked was the set of shfted deltas

$$\delta(t - \tau), \forall \tau \in \mathbb{R}$$

If we only used the linearity property, we'd have to test, and record, the effect of $\mathscr{H}$ on every single delta.
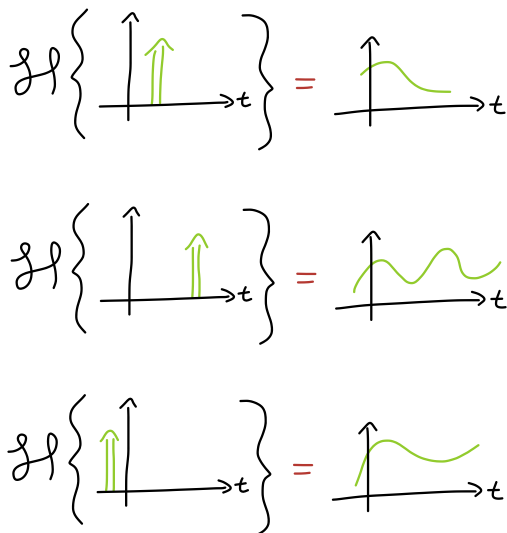
In other words, if we start by writing the input as a linear combination of the diracs (which is possible since they are a basis)

$$x(t) = \int_{-\infty}^{\infty} x(\tau)\delta(t - \tau)d\tau$$

And then apply $\mathscr{H}$ to this, we get[15]

$$\mathscr{H}\{x(t)\} = \mathscr{H}\left\{\int_{-\infty}^{\infty} x(\tau)\delta(t-\tau)d\tau\right\} = \left[\mathscr{H} \text{ is linear and } x(\tau) \text{ is constant wrt to t.}\right] =$$

$$= \int_{-\infty}^{\infty} x(\tau)\mathscr{H}\{\delta(t-\tau)\}d\tau$$

---

[15]**Be aware** that we're assuming that the system is *energy-free* in this derivation. We'll get to what that means later, but it turns out that assuming linearity bakes in energy-free:ness, so we're ok.

But now we use the time-invariance property

$$= \int_{-\infty}^{\infty} x(\tau) \mathscr{H}\{\delta(t)\}(t - \tau)d\tau$$

The amazing thing is that due to the time-invariance property, and the fact that the basis of dirac's are just shifted versions of each other, it suffices to just observe the effect of $\mathscr{H}\{\delta(t)\}$, to know the whole system!

Because of this special significance, we give $\mathscr{H}\{\delta(t)\}$ a special name; the "*impulse response*", and it's denoted by

$$\mathscr{H}\{\delta(t)\} = h(t)$$

The final integral is so commonly used that there's a special notation for it, namely

$$\int_{-\infty}^{\infty} x(\tau)h(t - \tau)d\tau = x(t) * h(t)$$

and is often called the *convolution* of x and h.

14.2   Frequency response

So we know plugging in a dirac into a system can yeild interesting results. In particular, since the shifted diracs are a basis then any (LTI[16]) system can be totally determined by its effect on it, which is why $\mathscr{H}\{\delta(t)\} = h(t)$, where $\mathscr{H}$ is the system, was significant.

What if we plug in a complex exponential into our (LTI) system?

$$\mathscr{H}\{e^{j\omega t}\} = e^{j\omega t} * h(t) = \int_{-\infty}^{\infty} e^{j\omega(t-\tau)}h(\tau)d\tau = e^{j\omega t}\left[\int_{-\infty}^{\infty} h(\tau)e^{-j\omega\tau}\right]$$

But notice now that $\int_{-\infty}^{\infty} h(\tau)e^{-j\omega\tau} = \mathscr{F}\{h(t)\}$. We'll denote this $H(\omega)$. Thus we have

$$\mathscr{H}\{e^{j\omega t}\} = H(\omega)e^{j\omega t}$$

Which means $e^{j\omega t}$ is an eigenfunction of LTI systems! Not only that, but these also make up a basis[17] (as shown[18] above), which means every system can be determined by its effect on the complex exponentials!

We often call $H(\omega)$ the frequency response.

14.3   Frequency response (alternative perspective)

If you don't care about the spectral theorem (which is what the above used to motivate the frequency response), check this out.

We already know that convolution in the time-domain becomes multiplication in the frequency-domain (just check out the "formelsamling"). That means that if we take the fourier transform of $y(t) = x(t) * h(t)$ and analyze it in the frequency spectrum, we get

$$Y(\omega) = X(\omega)H(\omega)$$

In other words, all an LTI system does is just increse of dampen the magnitude of the fourier transform of the input! The fourier spectrum of the impulse response, aka the frequency response, determines this increase/dampening effect.

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[16]Though it doesn't actually need to be time-invariant to be determined by the effect on the diracs.

[17]*Almost* orthonormal.

[18]Which was never really shown.

15 Inferring causality from $h(t)$

Another cool thing about the fact that $h(t)$ totally determines the system is that we can infer various properties of the system just by looking at $h(t)$. One of these things is causality.

So that a system is causal just means the output only depends on the past and current inputs, but not on future inputs. This can be seen from $h(t)$. A system is causal if and only if $h(t) = 0$, for all $t < 0$.
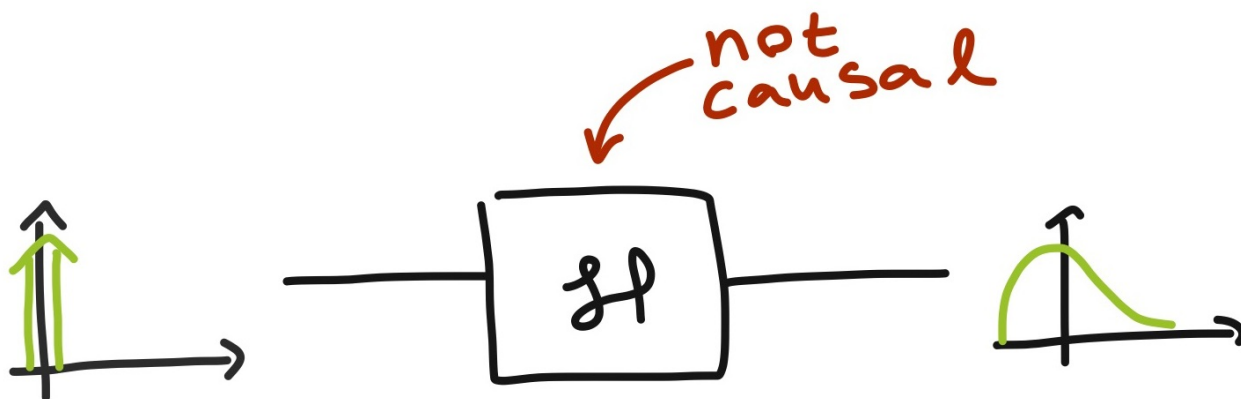
To motivate this, recall that $\mathscr{H}\{\delta(t)\} = h(t)$, thus if $h(t)$ does have a non-zero value for $t < 0$, that would mean when we sent an impulse into our system at $t = 0$, then the system would have already responded, e.g., at $t = -3$. In other words, the output would be able to predict the future.

16 Differential equations

If you've done any work with electronics, you know that a lot of systems can be described by a differential equation, i.e.

$$Q(D)y(t) = P(D)x(t)$$

where $Q(D), P(D)$ are polynomials of the differential operator, and $y(t)$ is the output, and $x(t)$ is the input. Plug in the input and you can solve for the output.

16.1 Energy

Before we discuss how to solve such a system we need to discuss what "energy" means in terms of systems. By definition, the energy $y_0(t)$, of the system $y(t) = \mathscr{H}\{x(t)\}$ is the output of the system when $x(t) = 0$, i.e. $y_0(t) = \mathscr{H}\{0\}$. That a system is *energy-free* means that $y_0(t) = 0$.

But if we assume the system is linear, then it's possible to show that $\mathscr{H}\{0\} = 0$. Check this out:

$$\mathscr{H}\{0\} = \mathscr{H}\{0 \cdot x(t)\} = \left[\text{ homogeneity }\right] = 0 \cdot \mathscr{H}\{x(t)\} = 0$$

So you'd think that if the system is LTI, then it's always energy-free, right? Well, not according to the lecturer. Because apparently the **L** in LTI can also mean "incrementally linear", which means "linear, but with $\mathscr{H}\{0\} \neq 0$"[19]. Thus if we're told our system is LTI, then your next question should always be "is it energy-free or not?". If the answer is "energy-free" then the **L** means "linear", if the answer if "not energy-free", then **L** means "incrementally linear".

In practice (on the exam), the energy will either be given explicitly or you'll be able to find it from the differential equation description of the system.

16.2   Solving the differential equation, and the transfer function

So given our system described by the differential equation

$$Q(D)y(t) = P(D)x(t)$$

how do we go about solving it?

We could just use our normal approach of finding the homogeneous and particular solution solutions. I'm not going to go into how we do that, but instead I'll just present a much simpler way involving the laplace transform.

First we'll need to assume our system is causal, i.e. $h(t) = 0$ for all $t < 0$, and that the input has the property that $x(t) = 0$ for all $t < 0$. Now let's expand the polynomials into

$$\left(\alpha_N \frac{d^N}{dt^N} + \alpha_{N-1}\frac{d^{N-1}}{dt^{N-1}} + \cdots + \alpha_1 \frac{d}{dt} + \alpha_0\right) y(t) = \left(\beta_M \frac{d^M}{dt^M} + \beta_{M-1}\frac{d^{M-1}}{dt^{M-1}} + \cdots + \beta_1 \frac{d}{dt} + \beta_0\right) x(t)$$

Now we use the formulas[20] on page 18, which states

$$\mathscr{L}_I\left\{\frac{dx(t)}{dt}\right\} = sX(s) - x(0^-)$$

and

$$\mathscr{L}_I\left\{\frac{d^2x(t)}{dt^2}\right\} = s^2 X(s) - sx(0^-) - x'(0^-)$$

and in general

$$\mathscr{L}_I\left\{\frac{d^n x(t)}{dt^n}\right\} = s^n X(s) - \sum_{k=1}^{n} s^{n-k}\frac{d^{k-1}x(t)}{dt^{k-1}}\bigg|_{t=0^-}$$

So if we take the unilateral laplace transform of the differential equation we get

$$\left(\alpha_N s^N + \cdots + \alpha_1 s + \alpha_0\right) Y(s) - Y_i(s) = \left(\beta_M s^M + \cdots + \beta_1 s + \beta_0\right) X(s)$$

where $Y_i(s)$ are all those $y(0^-), sy(0^-)-y'(0^-), \ldots$ terms. Note that the reason we don't get any $x(0^-), x'(0^-), \ldots$ terms is because we assumed that $x(t) = 0$ for all $t < 0$.

---

[19]I know, it sounds like a contradiction. In truth I'm not sure how "incrementally linear" could be defined more formally.
[20]I don't like to just use formulas, but I have no idea how to derive this.

Rewriting this a bit we get

$$Y(s) = Y_0(s) + \frac{\left(\beta_M s^M + \cdots + \beta_1 s + \beta_0\right)}{\left(\alpha_N s^N + \cdots + \alpha_1 s + \alpha_0\right)} X(s)$$

where $Y_0(s) = \frac{Y_i(s)}{(\alpha_N s^N + \cdots + \alpha_1 s + \alpha_0)}$. (Notice that we divide by $Q(s)$ without regard for whether or not it is ever zero. This is ISY math, not MAI math.)

Let's assume, for a moment, that $Y_0(s) = 0$, i.e. $y(0^-) = y'(0^-) = \cdots = \left.\frac{d^N y(t)}{dt^N}\right|_{t=0^-} = 0$ thus

$$Y(s) = H(s)X(s)$$

where $H(s) = \frac{\left(\beta_M s^M + \cdots + \beta_1 s + \beta_0\right)}{\left(\alpha_N s^N + \cdots + \alpha_1 s + \alpha_0\right)}$.

But doesn't that look suspiciously similar to something else we've seen? Namely that

$$Y(\omega) = H(\omega)X(\omega)$$

And it is for this reason that we call $H(s) = \frac{Q(s)}{P(s)}$ (assuming energy-free:ness), the "transfer function".

Except for our use of notation, we haven't really motivated this[21], but it turns out that not only is it true (by definition) that $H(s) = \frac{Q(s)}{P(s)}$, but it's also[22] true that $H(s) = \mathscr{L}_I\{h(t)\}$. which means that we can find the frequency response $H(\omega)$ by looking at the $j\omega$ axis of $H(s)$.

16.3   Inferring the frequency response from the poles/zeroes

So we know that if we know $H(s)$, then we can just plug in $s = j\omega$ to get the $H(\omega)$, which determines the entire system. This is great because it gives us a tool for intuitively and quickly reason about a system by looking at the poles and zeroes of the transfer function.

Suppose we plotted the zeroes and poles of $H(s)$ like so
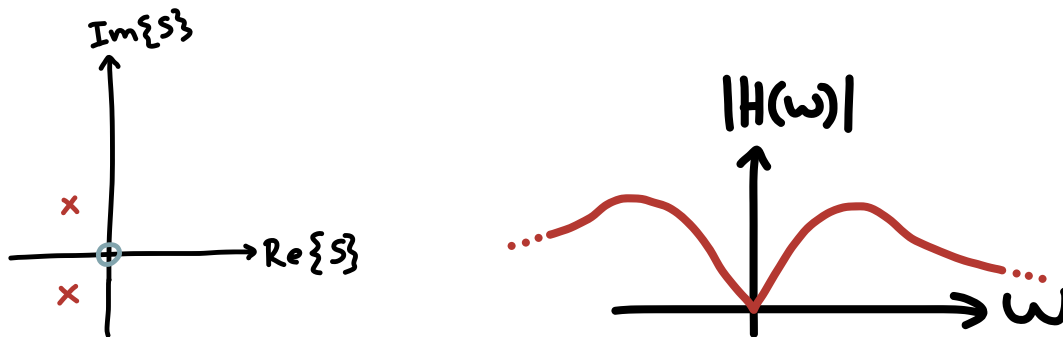
Since poles indicate where the $|H(s)|$ explode to infinity, and poles where $H(s) = 0$ we can sketch $H(\omega)$ purely from this illustration (assuming, of course, that $Re\{s\} = 0$ is in the ROC). This method becomes invaluable when designing LTI systems (aka filters).

---

[21]Neither does the lecurer (which is why I can't).
[22]I wish I could find a proof/motivation for this.

17 Stability

The properties of systems we've dicussed so far have been time-invariance, linearity, energy-free:ness and causality. Now it's time for stability. Before we even begin the discussion, let's assume our system $\mathscr{H}$ is LTI and causal.

Our system can either be internally stable or externally stable.

17.1 Internal stability

That a system is internally stable means that

$$\lim_{t \to \infty} y_0(t) = 0$$

This basically means that if we wait long enough, the energy won't be visible in the output anymore, which means that if we know our system is internally stable, then we can just wait for a few seconds (or minutes) and it'll then act truly linear (not incrementally linear). For this reason internal stability is also sometimes called *asymptotically stable*.

If you recall from the derivation above,

$$Y_0(s) = \frac{Y_i(s)}{(\alpha_N s^N + \cdots + \alpha_1 s + \alpha_0)}$$

and if you trust the lecturer[23] that $Y_0(s) = \mathscr{L}_I\{y_0(t)\}$, then you see that if the roots of the denominator have roots which all satisfy $Re\{s\} < 0$, then the system is internally stable (since you can do PBU on $Y_0(s)$, which means $y_0(t)$ is a bunch of exponentials whose arguments are the roots of polynomial). Apparently this is also true in the reverse direction; i.e. $\mathscr{H}$ is internally stable if and only if $Q(\lambda)$ has roots all satisfying $Re\{\lambda\} < 0$.

If all the roots of $Q(s)$ satisfy $Re\{s\} < 0$ then (also using PBU, and assuming the degree of $Q(D)$, aka N, is greater then or equal to M, the degree of $P(D)$) $\int_{-\infty}^{\infty} h(t)dt < \infty$, which implies that the system is externally stable, as we'll see below.

17.2 External stability

When we think about the external stability of the system we only care about the input and output of the system, not so much about what goes on inside (not so much about the differential equation). For this reason external stability is also known as *BIBO stability* (bounded-input-bounded-output). So for a system to be BIBO stable, all inputs $x(t)$ which satisfy

$$|x(t)| < K_1 < \infty$$

for some K_1, must produce an output $y(t)$ which is also bounded

$$|y(t)| < K_2 < \infty$$

for some K_2.

---

[23]Which isn't motivated in any way except by using clever tricks of notation.

"Every possible input"? This seems like an annoying thing to check. Luckily $h(t)$ comes to the rescue. First assume that $h(t)$ is absolutely integrable[24], i.e.

$$\int_{-\infty}^{\infty} |h(t)|dt = C < \infty$$

and that the input is bounded

$$|x(t)| < K < \infty$$

Now look what happends to the output

$$y(t) = h(t) * x(t) = \int_{-\infty}^{\infty} h(\tau)x(t-\tau)d\tau \leq \int_{-\infty}^{\infty} |h(\tau)||x(t-\tau)|d\tau \leq K \int_{-\infty}^{\infty} |h(\tau)|d\tau < K \cdot C$$

Thus $h(t)$ being absolutely integrable implies that the system is BIBO stable!

17.3   Marginal stability

Finally there's marginal stability. That a system is *marginally stable* means that

$$\lim_{t \to \infty} y_0(t) \notin \{0, \infty\}$$

This just means that at least one pole is on the $j\omega$ axis[25], and that the rest are on the left side of the $j\omega$ axis.

The fact that these statements are equivalent can also be motivated using PBU and the fact that the roots of $Q(\lambda)$ become the arguments of the exponentials $e^{\lambda t}$ in $h(t)$.

---

[24]That a function is absolutely integrable is sometimes given as

$$\int_{-\infty}^{\infty} |h(t)|^2 dt < \infty$$

but of course this is equivalent to
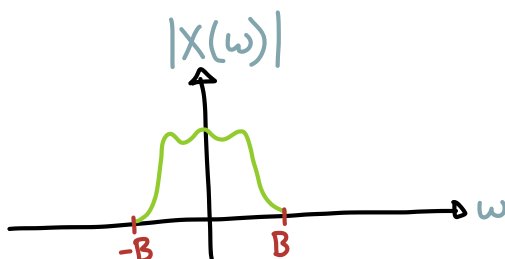
$$\int_{-\infty}^{\infty} |h(t)|dt < \infty$$

[25]There can't be any poles on the same place on the $j\omega$ axis though.

18   Modulation

Modulation is a way of moving a signal from one frequency band to another (usually higher). This is often a required operation in communication.

So suppose you have a signal $x(t)$ whose frequency-spectrum $X(\omega)$ looks like this.

And you want it to be centered at ω_0 instead. What do we always do when we want to move a signal? Convolve with a shifted dirac of course! So our modulated signal X_{ω_0} becomes

$$X_{\omega_0}(\omega) = X(\omega) * \delta(\omega - \omega_0)$$

Not so fast. Notice that this makes the frequency-spectrum asymetric, which means the corresponding signal in the time-domain will be complex. In the end we can only deal with signals in the time-domain. We can't send a complex-valued function through a speaker. To fix this we just convolve with a delta on both sides of the spectrum, i.e.
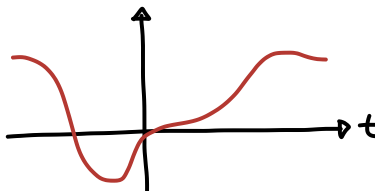
$$X_{\omega_0}(\omega) = X(\omega) * (\delta(\omega + \omega_0) + \delta(\omega - \omega_0))$$
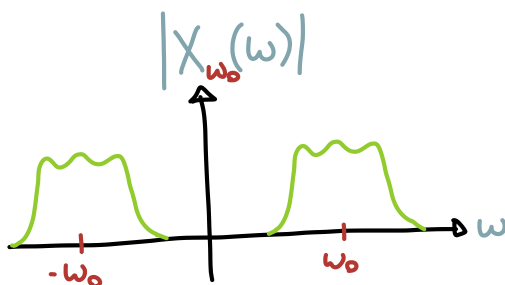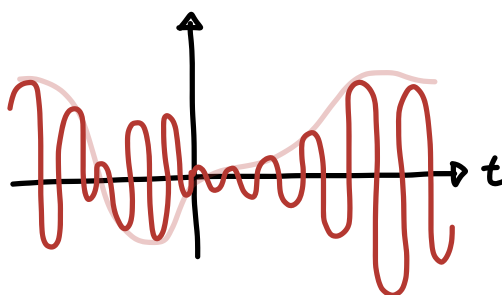
That's all well and good. What happends in the time-domain after this operation?

$$x_{\omega_0}(t) = \mathscr{F}^{-1}\{X_{\omega_0}\} = \mathscr{F}\{X(\omega) * (\delta(\omega + \omega_0) + \delta(\omega - \omega_0))\} = \frac{1}{\pi}x(t)cos(\omega_0 t)$$

In other words, if our original signal was


then our modulated signal would look like

18.1   Modulation is not LTI

Recall that if $H(\omega)$ is the frequecy response to an LTI system, then the output $Y(\omega)$ can be determined by $Y(\omega) = H(\omega)X(\omega)$, where $X(\omega)$ is the input. Notice thus, that LTI systems can only amplify or dampen existing frequecies of $X(\omega)$, but never introduce new frequencies (if $X(\omega_0) = 0$ then multiplying by $H(\omega_0)$ cannot produce anything higher than zero).

The whole point of modulation is to kill of frequencies at $\omega = 0$ and move them to some other place $\omega = \omega_0$. Thus modulation is not LTI.

18.2   Demodulation

So after we've modulated a signal and sent it over the air, how do we go about getting the original signal back? Suppose we got the following signal $X(\omega)$
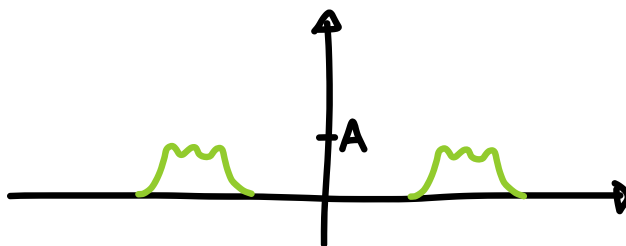
The way to demodulate this is to convolve with precisely the same diracs as before.
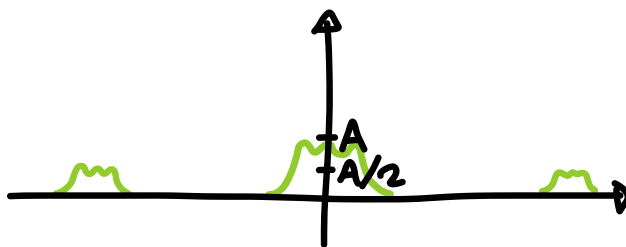
Notice that this produces some "*noise*" at the edges of the spectrum. To reduce this we can pass the spectrum through a low-pass filter.
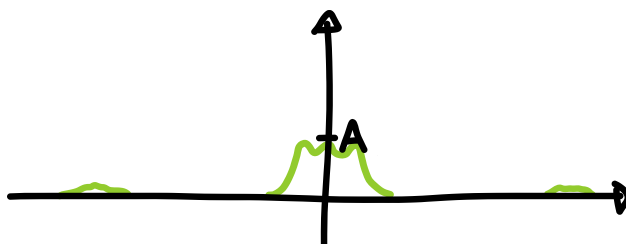
Realistic low-pass filters are rarely perfect, so some of that noise is usually still there. But we can at least reduce it.

19   Sampling

Now that we've spent the majority of our time in continous systems, it's finally time to move into the discrete world. To do that we must first return to the issue of sampling.
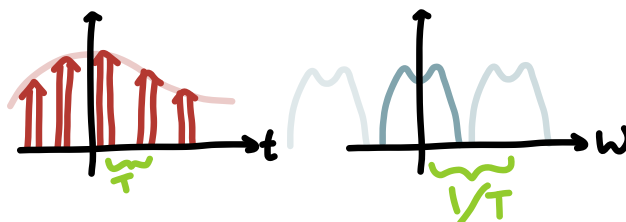
Sampling (ideally) is just multiplication with III_T;

$$\overline{x}(t) = \text{III}_T x(t)$$

But what happends in the fourier domain for the sampled signal? Well multiplication turns into convolution, so[26]

$$\overline{X}(\omega) = \frac{1}{T}\text{III}_{\frac{1}{T}} X(\omega)$$

Which means that the sampled signal have the frequecy spectrum of the original signal, but periodized at a frequency $f_s = \frac{1}{T}$.


---

[26]Why isn't the $\frac{1}{2\pi}$ factor here? It's there in the "formelsamling", but not in chapter 8 of the slides.

But at which frequency should you sample? Well what the sample-theorem tells us is that if we assume our signal $x(t)$ is band-limited, i.e. there exists an B such that $X(f) = 0$ for all $f > B$, then we should just sample at twice that frequency.

$$f_s = 2B$$

Why twice? Well, if we sample below that then the different shifted versions of the original frequency spectrums will start to overlap, which means we won't be able to reconstruct the signal back to its original form.
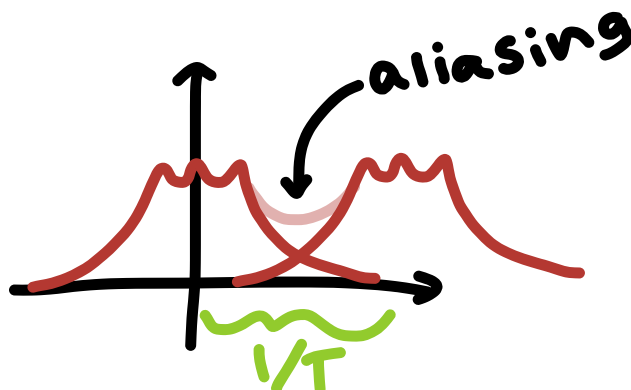
This effect is called aliasing. The practical effect is that when you under-sample a signal (e.g. a piece of music), then the higher frequencies will be more distorted then the lower frequencies.

20   Discrete

I'm not going to write much about the discrete world, because it's pretty much identical to the continuous world. Here are some of the key differences.

- Instead of continuous signals $x(t)$, we have discrete ones $x[n]$.

- Instead of the differential operator D, we use the "advancement operator" $Ex[n] = x[n + 1]$, and our differential equations turn into *difference equations*

$$\left(E^N + \alpha_{N-1}E^{N-1} + \cdots + \alpha_1 E^1 + \alpha_0\right) y[n] = \left(E^N + \beta_{N-1}E^{N-1} + \cdots + \beta_1 E^1 + \beta_0\right) x[n]$$

and all the stuff about energy-free:ness, stability, etc are the same/similar.

- Instead of the laplace transform we use the z-transform.

21 Things I don't know

There are lots of things I don't understand or have forgotten. Here's a list of known unknows[27].

- Why is $\mathscr{L}_I\left\{\frac{dx(t)}{dt}\right\} = sX(s) - x(0^-)$?

- When solving the differential equation using the laplace transform, we had to assume causality. What do we do if the system isn't causal?

- Why can't we just plug in a dirac directly into $Q(D)y(t) = P(D)x(t)$, then solve for $y(t)$ and take the laplace transform?

- I'm not totally sure about how to derive (i.e. motivate) the $T_0 = N_0 T$ and $f_s = N_0 f_0$ thing for the DFT, which is why I didn't write anything about it.

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[27]Of course, there are probably plenty of unknown unknows too.