# Lecture notes 4: Fourier Analysis

#### **Definitions**

There are many common (and confusing, but ultimately trivial!) differences in defining the *Fourier transform*. One common definition is

$$F(\nu) = \int_{-\infty}^{\infty} f(t)e^{-i2\pi\nu t}dt$$

Thus  $F(\nu)$  gives the wavenumber representation of the function f(t). The inverse transform can be written

$$f(t) = \int_{-\infty}^{\infty} F(\nu)e^{i2\pi\nu t}d\nu$$

 $F(\nu)$  is in general a complex function whose interpretation may be aided by expression in the polar coordinate form  $F(\nu) = A(\nu)e^{i\phi(\nu)}$ , where  $A(\nu)$  and  $\phi(\nu)$  are real functions where  $A(\nu) = |F(\nu)|$  is the amplitude and  $\phi(\nu) = \arg[F(\nu)]$  is the phase. Note that we then can write the inverse transform as

$$f(t) = \int_{-\infty}^{\infty} A(\nu)e^{i[2\pi\nu t + \phi(\nu)]}d\nu,$$

which is seen to be a recombination of all the frequency components of f(t). Each component is a complex sinusoid of the form  $e^{2\pi i\nu t}$  whose amplitude is  $A(\nu)$  and whose initial phase (at t=0) is  $\phi(\nu)$ . This interpretation of the Fourier transform clearly shows its relation to the Fourier series<sup>1</sup>

$$f(x) = \frac{a_0}{2} + \sum_{n=1}^{\infty} [a_n \cos(nx) + b_n \sin(nx)]$$

with coefficients given by

$$a_n = \frac{1}{\pi} \int_{-\pi}^{\pi} f(x) \cos(nx) dx$$
$$b_n = \frac{1}{\pi} \int_{-\pi}^{\pi} f(x) \sin(nx) dx.$$

It is common to perform the substitution  $\nu = \omega/2\pi$  which gives

$$\begin{split} F(\omega) &= \int_{-\infty}^{\infty} f(t) e^{i\omega t} dt \\ f(t) &= \frac{1}{2\pi} \int_{-\infty}^{\infty} F(\omega) e^{-i\omega t} d\omega \end{split}$$

<sup>&</sup>lt;sup>1</sup>Note that by using Euler's formula  $e^{inx} = \cos(nx) + i\sin(nx)$  a more concise (and modern) form can be used:  $f(x) = \sum_{n=-\infty}^{\infty} c_n e^{inx}$  with  $c_n = \frac{1}{2\pi} \int_{-\pi}^{\pi} f(x) e^{-inx} dx$ .

## **Properties**

Fourier transforms exhibit a number of properties that are very useful:

• Linearity

$$af(t) + bg(t) \Longleftrightarrow aF(\omega) + bG(\omega)$$

• Multiplication

$$f(t)g(t) \Longleftrightarrow \frac{1}{2\pi}(F \otimes G)(\omega)$$

where we define the  $convolution \otimes$  as

$$K(\omega) \equiv \int_{-\infty}^{\infty} F(\omega') G(\omega - \omega') d\omega'$$

• Convolution

$$(f \otimes g)(t) \Longleftrightarrow F(\omega)G(\omega)$$

• Conjugation

$$\overline{f(t)} \Longleftrightarrow \overline{F(-\omega)}$$

• Scaling

$$f(at) \Longleftrightarrow \frac{1}{|a|} F\left(\frac{\omega}{a}\right)$$

• Time reversal

$$f(-t) \iff F(-\omega)$$

• Time shift

$$f(t-t_0) \Longleftrightarrow e^{-i\omega t_0} F(\omega)$$

• Parsevals theorem

$$\int_{-\infty}^{\infty} f(t) \overline{g(t)} dt = \frac{1}{2\pi} \int_{-\infty}^{\infty} F(\omega) \overline{G(\omega)} d\omega$$

The correlation of two functions, denoted by Corr(g, h), is defined

$$Corr(g,h) \equiv \int_{-\infty}^{\infty} g(t'+t)h(t')dt'$$

The correlation is a function of t, called the lag. The "Correlation theorem" states that

$$\operatorname{Corr}(g,h) \Longleftrightarrow G(f)\overline{H(f)}$$

if g and h are real functions, as we expect for our applications [otherwise the expression on the right hand side of the correlation theorem is G(f)H(-f).] The correlation of a function with itself is called the *autocorrelation*. In this case we have

$$\operatorname{Corr}(g,g) \Longleftrightarrow |G(f)|^2$$

which is also known as the "Wiener-Khinchin theorem". The *total power* in a signal can be computed from the autocorrelation of that signal and following Parseval's theorem we can write

Total Power 
$$\equiv \int_{-\infty}^{\infty} |g(t)|^2 dt = \int_{-\infty}^{\infty} |G(f)|^2 df$$
.

Often one desires to know how much power is contained in the frequency interval between f and f + df. In this case it is often not interesting to distinguish between positive and negative f, but rather regard f as varying from 0 (zero frequency or DC) to  $+\infty$ . In this case we define the *one-sided power spectral density* of the function g:

$$P_q(f) \equiv |G(f)|^2 + |G(-f)|^2 \qquad 0 \le f < \infty,$$

when q is real the two terms in the equation above are equal, so

$$P_g(f) = 2 |G(f)|^2$$
.

## Sampling Theorem and Aliasing

We will in general be dealing with functions h(t) which are sampled at evenly spaced intervals in time (or space). Let  $\Delta$  denote the time (space) interval between consecutive samples.

For any sampling interval  $\Delta$ , there is a special frequency  $\nu_c$  called the *Nyquist* frequency given by

$$\nu_c \equiv \frac{1}{2\Delta}$$

The critical sampling of a sine wave is two sample points per cycle. There are two aspects of the critical frequency. First, if the original signal is bandwidth limited to frequencies smaller than  $\nu_c$  then the function is completely determined by its discrete samples. This is the *sampling theorem*. However, if a signal is *not* bandwidth limited to less than the Nyquist frequency then the power that lies outside the range  $-\nu_c < \nu < \nu_c$  is spuriously moved into that range. This phenomena is called *aliasing*.

#### **FFTs**

Let us now estimate a Fourier transform from a finite number of sampled points. Suppose that we have N consecutive sampled values

$$h_k \equiv h(t_k)$$
  $t_k \equiv k\Delta$   $k = 0, 1, 2, \dots, N-1$ 

so that the sampling interval is  $\Delta$ . Also assume N is even. With N numbers of input, we cannot produce more than N independent numbers of output. So, instead of trying to estimate the Fourier transform  $H(\nu)$  at all values of  $\nu$  in the range  $-\nu_c$  to  $\nu_c$ , let us seek estimates at only the discrete values

$$\nu_n \equiv \frac{n}{N\Lambda}, \qquad n = \frac{-N}{2}, \dots, \frac{N}{2}$$

The extreme values of n correspond to the lower and upper values of the Nyquist critical frequency range.

The remaining step is to approximate the continuous transform by a discrete sum

$$H(\nu_n) = \int_{-\infty}^{\infty} h(t)e^{2\pi i\nu_n t}dt \approx \sum_{k=0}^{N-1} h_k e^{2\pi i\nu_n t_k} \Delta = \Delta \sum_{k=0}^{N-1} h_k e^{2\pi ikn/N}$$

The final summation is called the discrete Fourier transform of the N points  $h_k$ . Let us denote it by  $H_n$ :

$$H_n \equiv \sum_{k=0}^{N-1} h_k e^{2\pi i k n/N} \tag{1}$$

Up to now we have taken the view that index n varies from -N/2 to N/2. However, since it is clear that equation 1 is periodic in n we also have that  $H_{-n}=H_{N-1}$ . With this in mind it is customary to let n vary from 0 to N-1 (one period). When this convention is followed the zero frequency corresponds to n=0, positive frequencies  $0<\nu<\nu_c$  correspond to values  $1\leq n\leq N/2-1$ , while negative frequencies  $-\nu_c<\nu<0$  correspond to  $N/2_1\leq n\leq N-1$ . The value n=N/2 corresponds to both  $\nu=\nu_c$  and  $\nu=-\nu_c$ .

The discrete *inverse* transform which recovers  $h_k$  form the  $H_n$  is

$$h_k = \frac{1}{N} \sum_{n=0}^{N-1} H_n e^{-2\pi i k n/N}$$

How do we implement the discrete transform in code? The brute strength approach takes of order  $N^2$  operations: Define W as the complex number

$$W \equiv e^{2\pi i/N}$$

Then equation 1 can be written

$$H_n = \sum_{k=0}^{N-1} W^{nk} h_k,$$

i.e. the vector  $h_k$  is multiplied by a matrix whose  $(n,k)^{th}$  element is the constant W to the power  $n \times k$ . The matrix multiplication produces a vector result whose components are  $H_n$ . This multiplication evidently needs  $N^2$  complex multiplications.

In fact, the discrete Fourier transform can be achieved in  $N \log_2 N$  operations with an algorithm called the it Fast Fourier Transform or FFT. Here is one derivation of the FFT, that of Danielson and Lanczos in 1942. They showed that a discrete transform of length N can be rewritten as the sum of two discrete transforms, each of length N/2. One of the two is formed from the even

numbered points of the original N, the other from the odd-numbered points.

$$F_{k} = \sum_{j=0}^{N-1} e^{2\pi i j k/N} f_{j}$$

$$= \sum_{j=0}^{N/2-1} e^{2\pi i 2 j k/N} f_{2j} + \sum_{j=0}^{N/2-1} e^{2\pi i (2j+1)k/N} f_{2j+1}$$

$$= \sum_{j=0}^{N/2-1} e^{2\pi i j k/(N/2)} f_{2j} + W^{k} \sum_{j=0}^{N/2-1} e^{2\pi i j k/(N/2)} f_{2j+1}$$

$$= F_{k}^{e} + W^{k} F_{k}^{o}$$

 $F_k^e$  denotes the  $k^{th}$  component of the Fourier transform of length (N/2) formed from the even components of the original  $f_j$ , while  $F_k^o$  is the corresponding transform of length (N/2) from the odd components.

This procedure can be applied recursively; having reduced the problem of computing  $F_k$  to that of computing  $F_k^e$  and  $F_k^o$ , one can do the same reduction of  $F_k^e$  to the problem of computing the transform of its N/4 even-numbered inputs data,  $F_k^{ee}$ , and N/4 odd-numbered data  $F_k^{eo}$ . When we start with an original N which is a integer power of 2 (one should only use FFTs with N a power of 2, padding the input data with zeroes is necessary) we can continue applying the Danielson-Lanczos method until the original data is subdivided all the way down to transforms of length one. The Fourier transform of length one is just the identity operation that copies its one input number into its output slot. Thus, for every pattern of e's and e's numbering  $\log_2 N$  in all there is a one-point transform that is given by

$$F_k^{eoeeoeo\cdots oee} = f_n$$
 for some  $n$ 

Which value of n corresponds to which pattern? Reverse the pattern of e's and o's, let e=0 and o=1 and one will have, in binary, the value of n. This is because the successive subdivisions of the data into odd and even are tests of successive least significant bits of n. Thus by rearranging the input vector  $f_j$  in bit reversed order so that the individual numbers are in the order obtained by bit reversing j. The points are given as one-point transforms. These are recombined with the adjacent number to give two-point transforms, then combine adjacent pairs again to give 4-point transforms, and so on, using the Danielson-Lanczos formula at every step.

The FFT therefore has a structure where first the data are sorted into bit reversed order and thereafter the transforms of length  $2,4,8,\ldots,N$  are computed implementing the Danielson-Lanczos formula.

#### Exercises

Experiment with the fft function in idl. First make an x-axis, eg idl> x=findgen(2000)/100.\*2.\*!pi

- 1. Compute the amplitude of the transform of  $f = \sin x$ . In plotting this amplitude, what should be used for the x-axis?
- 2. What happens if you set the edges of f to zero: eg multiply f by a step-function s such as idl> s=fltarr(2000) & s[200:1800]=1.0? Overplot the amplitude of the transform as you narrow the region of s that is equal to one.
- 3. Compute the transform of step-functions of various widths.
- 4. Add sinusoidal functions to f with different frequencies and check the resulting transforms. What happens to the transform if you add a constant to f?
- 5. Consider a function g given by the sum of three sinusoidals of differing frequency. Construct such a g and remove one of the frequencies from g using forward and back transforms FFT's in idl.