



WIKIPEDIA
The Free Encyclopedia

Main page
Contents
Featured content
Current events
Random article
Donate to Wikipedia
Wikipedia store

Interaction

Help
About Wikipedia
Community portal
Recent changes
Contact page

Tools

What links here
Related changes
Upload file
Special pages
Permanent link
Page information
Wikidata item
Cite this page

Print/export

Create a book
Download as PDF
Printable version

Languages

العربية
Català
Čeština
Deutsch
Español
فارسی
Français
한국어
हिन्दी
Bahasa Indonesia
Italiano
Lietuvių
Nederlands
日本語
Polski
Português
Русский
Српски / srpski
Srpskohrvatski /
српскохрватски
Basa Sunda
Suomi
Svenska
Türkçe
Українська
Tiếng Việt
中文

Edit links

Create account Log in

Article Talk

Read Edit View history

Search



Discrete Fourier transform

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In **mathematics**, the **discrete Fourier transform (DFT)** converts a finite list of equally spaced **samples** of a function into the list of **coefficients** of a finite combination of **complex sinusoids**, ordered by their **frequencies**, that has those same sample values. It can be said to convert the sampled function from its original domain (often **time** or position along a line) to the **frequency domain**.

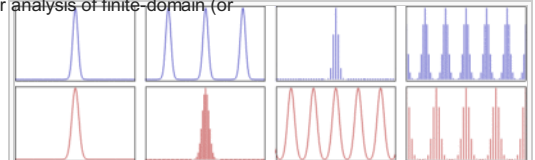
The input samples are **complex numbers** (in practice, usually **real numbers**), and the output coefficients are complex as well. The frequencies of the output sinusoids are integer multiples of a fundamental frequency, whose corresponding period is the length of the sampling interval. The combination of sinusoids obtained through the DFT is therefore **periodic** with that same period. The DFT differs from the **discrete-time Fourier transform (DTFT)** in that its input **and** output sequences are both finite; it is therefore said to be the Fourier **analysis of finite-domain** (or periodic) discrete-time functions.

The DFT is the most important **discrete transform**, used to perform **Fourier analysis** in many practical applications.^[1] In **digital signal processing**, the function is any quantity or **signal** that varies over time, such as the pressure of a **sound wave**, a **radio** signal, or daily **temperature** readings, sampled over a finite time interval (often defined by a **window function**^[2]). In **image processing**, the samples can be the values of **pixels** along a row or column of a **raster image**. The DFT is also used to efficiently solve **partial differential equations**, and to perform other operations such as **convolutions** or multiplying large integers.

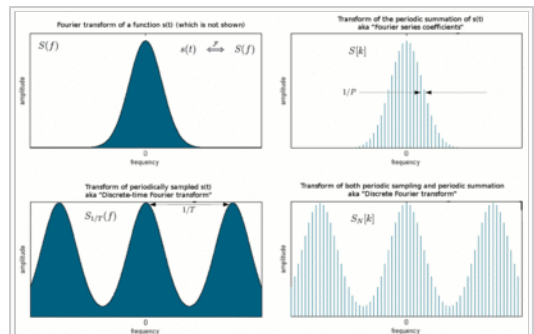
Since it deals with a finite amount of data, it can be implemented in **computers** by **numerical algorithms** or even dedicated **hardware**. These implementations usually employ efficient **fast Fourier transform (FFT)** algorithms;^[3] so much so that the terms "FFT" and "DFT" are often used interchangeably. Prior to its current usage, the "FFT" **initialism** may have also been used for the ambiguous term "finite Fourier transform".

Fourier transforms

Continuous Fourier transform
Fourier series
Discrete-time Fourier transform
Discrete Fourier transform
Fourier analysis
Related transforms



Relationship between the (continuous) **Fourier transform** and the discrete Fourier transform. **Left column:** A continuous function (top) and its Fourier transform (bottom). **Center-left column:** **Periodic summation** of the original function (top). Fourier transform (bottom) is zero except at discrete points. The inverse transform is a sum of sinusoids called **Fourier series**. **Center-right column:** Original function is discretized (multiplied by a **Dirac comb**) (top). Its Fourier transform (bottom) is a periodic summation (**DTFT**) of the original transform. **Right column:** The DFT (bottom) computes discrete samples of the continuous DTFT. The inverse DFT (top) is a periodic summation of the original samples. The **FFT** algorithm computes one cycle of the DFT and its inverse is one cycle of the DFT inverse.



Depiction of a Fourier transform (upper left) and its periodic summation (DTFT) in the lower left corner. The spectral sequences at (a) upper right and (b) lower right are respectively computed from (a) one cycle of the periodic summation of $s(t)$ and (b) one cycle of the periodic summation of the $s(nT)$ sequence. The respective formulas are (a) the **Fourier series integral** and (b) the **DFT summation**. Its similarities to the original transform, $S(f)$, and its relative computational ease are often the motivation for computing a DFT sequence.

Contents

[hide]

- Definition
- Properties
 - 2.1 Completeness
 - 2.2 Orthogonality
 - 2.3 The Plancherel theorem and Parseval's theorem
 - 2.4 Periodicity
 - 2.5 Shift theorem
 - 2.6 Circular convolution theorem and cross-correlation theorem
 - 2.7 Convolution theorem duality
 - 2.8 Trigonometric interpolation polynomial
 - 2.9 The unitary DFT
 - 2.10 Expressing the inverse DFT in terms of the DFT
 - 2.11 Eigenvalues and eigenvectors
 - 2.12 Uncertainty principle
 - 2.13 The real-input DFT
- Generalized DFT (shifted and non-linear phase)
- Multidimensional DFT
 - 4.1 The real-input multidimensional DFT
- Applications
 - 5.1 Spectral analysis
 - 5.2 Filter bank
 - 5.3 Data compression
 - 5.4 Partial differential equations
 - 5.5 Polynomial multiplication
 - 5.5.1 Multiplication of large integers
 - 5.5.2 Convolution
- Some discrete Fourier transform pairs
- Generalizations
 - 7.1 Representation theory
 - 7.2 Other fields
 - 7.3 Other finite groups
- Alternatives
- See also
- Notes
- Citations
- References
- External links

Definition

[edit]

The sequence of N complex numbers x_0, x_1, \dots, x_{N-1} is transformed into an N -periodic sequence of complex numbers:

$$X_k \stackrel{\text{def}}{=} \sum_{n=0}^{N-1} x_n \cdot e^{-j2\pi kn/N}, \quad k \in \mathbb{Z} \text{ (integers)} \quad [\text{note 1}] \quad (\text{Eq.1})$$

Because of [periodicity](#), the customary domain of k actually computed is $[0, N-1]$. That is always the case when the DFT is implemented via the [Fast Fourier transform](#) algorithm. But other common domains are $[-N/2, N/2-1]$ (N even) and $[-(N-1)/2, (N-1)/2]$ (N odd), as when the left and right halves of an FFT output sequence are swapped.^[4]

The transform is sometimes denoted by the symbol \mathcal{F} , as in $\mathbf{X} = \mathcal{F}\{\mathbf{x}\}$ or $\mathcal{F}(\mathbf{x})$ or $\mathcal{F}_\mathbf{x}$.^[note 2]

Eq.1 can be interpreted or derived in various ways, for example:

- It completely describes the [discrete-time Fourier transform](#) (DTFT) of an N -periodic sequence, which comprises only discrete frequency components. ([Using the DTFT with periodic data](#))
- It can also provide uniformly spaced samples of the continuous DTFT of a finite length sequence. ([Sampling the DTFT](#))
- It is the [cross correlation](#) of the *input* sequence, x_n , and a complex sinusoid at frequency k/N . Thus it acts like a [matched filter](#) for that frequency.
- It is the discrete analogy of the formula for the coefficients of a [Fourier series](#):

$$x_n = \frac{1}{N} \sum_{k=0}^{N-1} X_k \cdot e^{j2\pi kn/N}, \quad n \in \mathbb{Z} \quad (\text{Eq.2})$$

which is also N -periodic. In the domain $n \in [0, N-1]$, this is the **inverse transform** of [Eq.1](#). In this interpretation, each X_k is a complex number that encodes both amplitude and phase of a sinusoidal component ($e^{j2\pi kn/N}$) of function x_n . The sinusoid's [frequency](#) is k cycles per N samples. Its amplitude and phase are:

$$|X_k|/N = \sqrt{\text{Re}(X_k)^2 + \text{Im}(X_k)^2}/N$$

$$\arg(X_k) = \text{atan2}(\text{Im}(X_k), \text{Re}(X_k)) = -i \ln \left(\frac{X_k}{|X_k|} \right),$$

where [atan2](#) is the two-argument form of the [arctan](#) function.

The normalization factor multiplying the DFT and IDFT (here 1 and $1/N$) and the signs of the exponents are merely [conventions](#), and differ in some treatments. The only requirements of these conventions are that the DFT and IDFT have opposite-sign exponents and that the product of their normalization factors be $1/N$. A normalization of $\sqrt{1/N}$ for both the DFT and IDFT, for instance, makes the transforms unitary.

In the following discussion the terms "sequence" and "vector" will be considered interchangeable.

Using [Euler's Formula](#), it can be derived further to the forms commonly used in Engineering and Computer Science.

Fourier Transform:

$$X_k = \sum_{n=0}^{N-1} x_n \cdot \left(\cos(-2\pi k \frac{n}{N}) + j \sin(-2\pi k \frac{n}{N}) \right), \quad n \in \mathbb{Z} \quad (\text{Eq.3})$$

Inverse Fourier Transform:

$$x_n = \frac{1}{N} \sum_{k=0}^{N-1} X_k \cdot \left(\cos(2\pi k \frac{n}{N}) + j \sin(2\pi k \frac{n}{N}) \right), \quad n \in \mathbb{Z} \quad (\text{Eq.4})$$

N = number of time samples we have

n = current sample we're considering (0... $N-1$)

x_n = value of the signal at time n

k = current frequency we're considering (0 Hertz up to $N-1$ Hertz)

X_k = amount of frequency k in the signal (Amplitude and Phase, a complex number)

Properties [\[edit\]](#)

Completeness [\[edit\]](#)

The discrete Fourier transform is an invertible, [linear transformation](#)

$$\mathcal{F}: \mathbb{C}^N \rightarrow \mathbb{C}^N$$

with \mathbb{C} denoting the set of [complex numbers](#). In other words, for any $N > 0$, an N -dimensional complex vector has a DFT and an IDFT which are in turn N -dimensional complex vectors.

Orthogonality [\[edit\]](#)

The vectors $u_k = \left[e^{\frac{2\pi i}{N} kn} \mid n = 0, 1, \dots, N-1 \right]^T$ form an [orthogonal basis](#) over the set of N -dimensional complex vectors:

$$u_k^T u_{k'}^* = \sum_{n=0}^{N-1} \left(e^{\frac{2\pi i}{N} kn} \right) \left(e^{\frac{2\pi i}{N} (-k')n} \right) = \sum_{n=0}^{N-1} e^{\frac{2\pi i}{N} (k-k')n} = N \delta_{kk'}$$

where $\delta_{kk'}$ is the [Kronecker delta](#). (In the last step, the summation is trivial if $k = k'$, where it is $1+1+\dots=N$, and otherwise is a [geometric series](#) that can be explicitly summed to obtain zero.) This orthogonality condition can be used to derive the formula for the IDFT from the definition of the DFT, and is equivalent to the unitarity property below.

The Plancherel theorem and Parseval's theorem [\[edit\]](#)

If X_k and Y_k are the DFTs of x_n and y_n respectively then the [Parseval's theorem](#) states:

$$\sum_{n=0}^{N-1} x_n y_n^* = \frac{1}{N} \sum_{k=0}^{N-1} X_k Y_k^*$$

where the star denotes [complex conjugation](#). [Plancherel theorem](#) is a special case of the Parseval's theorem and states:

$$\sum_{n=0}^{N-1} |x_n|^2 = \frac{1}{N} \sum_{k=0}^{N-1} |X_k|^2.$$

These theorems are also equivalent to the unitary condition below.

Periodicity [\[edit\]](#)

The periodicity can be shown directly from the definition:

$$X_{k+N} \stackrel{\text{def}}{=} \sum_{n=0}^{N-1} x_n e^{-\frac{2\pi i}{N}(k+N)n} = \sum_{n=0}^{N-1} x_n e^{-\frac{2\pi i}{N}kn} \underbrace{e^{-2\pi i n}}_1 = \sum_{n=0}^{N-1} x_n e^{-\frac{2\pi i}{N}kn} = X_k.$$

Similarly, it can be shown that the IDFT formula leads to a periodic extension.

Shift theorem [\[edit\]](#)

Multiplying x_n by a *linear phase* $e^{\frac{2\pi i}{N}nm}$ for some integer m corresponds to a *circular shift* of the output X_k : X_k is replaced by X_{k-m} , where the subscript is interpreted [modulo](#) N (i.e., periodically). Similarly, a circular shift of the input x_n corresponds to multiplying the output X_k by a linear phase. Mathematically, if $\{x_n\}$ represents the vector \mathbf{x} then

$$\begin{aligned} \text{if } \mathcal{F}(\{x_n\})_k &= X_k \\ \text{then } \mathcal{F}(\{x_n \cdot e^{\frac{2\pi i}{N}nm}\})_k &= X_{k-m} \\ \text{and } \mathcal{F}(\{x_{n-m}\})_k &= X_k \cdot e^{-\frac{2\pi i}{N}km} \end{aligned}$$

Circular convolution theorem and cross-correlation theorem [\[edit\]](#)

The [convolution theorem](#) for the [discrete-time Fourier transform](#) indicates that a convolution of two infinite sequences can be obtained as the inverse transform of the product of the individual transforms. An important simplification occurs when the sequences are of finite length, \mathbf{N} . In terms of the DFT and inverse DFT, it can be written as follows:

$$\mathcal{F}^{-1}\{\mathbf{X} \cdot \mathbf{Y}\}_n = \sum_{l=0}^{N-1} x_l \cdot (y_N)_{n-l} \stackrel{\text{def}}{=} (\mathbf{x} * \mathbf{y}_N)_n,$$

which is the convolution of the \mathbf{x} sequence with a \mathbf{y} sequence extended by [periodic summation](#):

$$(\mathbf{y}_N)_n \stackrel{\text{def}}{=} \sum_{p=-\infty}^{\infty} y_{(n-pN)} = y_{(n \bmod N)}.$$

Similarly, the [cross-correlation](#) of \mathbf{x} and \mathbf{y}_N is given by:

$$\mathcal{F}^{-1}\{\mathbf{X}^* \cdot \mathbf{Y}\}_n = \sum_{l=0}^{N-1} x_l^* \cdot (y_N)_{n+l} \stackrel{\text{def}}{=} (\mathbf{x} \star \mathbf{y}_N)_n.$$

When either sequence contains a string of zeros, of length L , $L+1$ of the circular convolution outputs are equivalent to values of $\mathbf{x} * \mathbf{y}$. Methods have also been developed to use this property as part of an efficient process that constructs $\mathbf{x} * \mathbf{y}$ with an \mathbf{x} or \mathbf{y} sequence potentially much longer than the practical transform size (\mathbf{N}). Two such methods are called [overlap-save](#) and [overlap-add](#).^[5] The efficiency results from the fact that a direct evaluation of either summation (above) requires $O(N^2)$ operations for an output sequence of length N . An indirect method, using transforms, can take advantage of the $O(N \log N)$ efficiency of the [fast Fourier transform](#) (FFT) to achieve much better performance. Furthermore, convolutions can be used to efficiently compute DFTs via [Rader's FFT algorithm](#) and [Bluestein's FFT algorithm](#).

Convolution theorem duality [\[edit\]](#)

It can also be shown that:

$$\begin{aligned} \mathcal{F}\{\mathbf{x} \cdot \mathbf{y}\}_k &\stackrel{\text{def}}{=} \sum_{n=0}^{N-1} x_n \cdot y_n \cdot e^{-\frac{2\pi i}{N}kn} \\ &= \frac{1}{N} (\mathbf{X} * \mathbf{Y}_N)_k, \text{ which is the circular convolution of } \mathbf{X} \text{ and } \mathbf{Y}. \end{aligned}$$

Trigonometric interpolation polynomial [\[edit\]](#)

The [trigonometric interpolation polynomial](#)

$$\begin{aligned} p(t) &= \frac{1}{N} \left[X_0 + X_1 e^{2\pi i t} + \dots + X_{N/2-1} e^{(N/2-1)2\pi i t} + X_{N/2} \cos(N\pi t) + X_{N/2+1} e^{(-N/2+1)2\pi i t} + \dots + X_{N-1} e^{-2\pi i t} \right] \\ &\text{for } N \text{ even,} \\ p(t) &= \frac{1}{N} \left[X_0 + X_1 e^{2\pi i t} + \dots + X_{\lfloor N/2 \rfloor} e^{\lfloor N/2 \rfloor 2\pi i t} + X_{\lfloor N/2 \rfloor + 1} e^{-\lfloor N/2 \rfloor 2\pi i t} + \dots + X_{N-1} e^{-2\pi i t} \right] \text{ for } N \text{ odd,} \end{aligned}$$

where the coefficients X_k are given by the DFT of x_n above, satisfies the interpolation property $p(n/N) = x_n$ for $n = 0, \dots, N-1$.

For even N , notice that the [Nyquist component](#) $\frac{X_{N/2}}{N} \cos(N\pi t)$ is handled specially.

This interpolation is *not unique*: aliasing implies that one could add N to any of the complex-sinusoid frequencies (e.g. changing e^{-it} to $e^{i(N-1)t}$) without changing the interpolation property, but giving *different* values in between the x_n points. The choice above, however, is typical because it has two useful properties. First, it consists of sinusoids whose frequencies have the smallest possible magnitudes: the interpolation is [bandlimited](#). Second, if the x_n are real numbers, then $p(t)$ is real as well.

In contrast, the most obvious trigonometric interpolation polynomial is the one in which the frequencies range from 0 to $N - 1$ (instead of roughly $-N/2$ to $+N/2$ as above), similar to the inverse DFT formula. This interpolation does *not* minimize the slope, and is *not* generally real-valued for real x_n ; its use is a common mistake.

The unitary DFT [\[edit\]](#)

Another way of looking at the DFT is to note that in the above discussion, the DFT can be expressed as a [Vandermonde matrix](#), introduced by [Sylvester](#) in 1867,

$$\mathbf{F} = \begin{bmatrix} \omega_N^{0 \cdot 0} & \omega_N^{0 \cdot 1} & \dots & \omega_N^{0 \cdot (N-1)} \\ \omega_N^{1 \cdot 0} & \omega_N^{1 \cdot 1} & \dots & \omega_N^{1 \cdot (N-1)} \\ \vdots & \vdots & \ddots & \vdots \\ \omega_N^{(N-1) \cdot 0} & \omega_N^{(N-1) \cdot 1} & \dots & \omega_N^{(N-1) \cdot (N-1)} \end{bmatrix}$$

where

$$\omega_N = e^{-2\pi i/N}$$

is a primitive [Nth root of unity](#).

The inverse transform is then given by the inverse of the above matrix,

$$\mathbf{F}^{-1} = \frac{1}{N} \mathbf{F}^*$$

With [unitary](#) normalization constants $1/\sqrt{N}$, the DFT becomes a [unitary transformation](#), defined by a unitary matrix:

$$\begin{aligned} \mathbf{U} &= \mathbf{F}/\sqrt{N} \\ \mathbf{U}^{-1} &= \mathbf{U}^* \\ |\det(\mathbf{U})| &= 1 \end{aligned}$$

where *det()* is the [determinant](#) function. The determinant is the product of the eigenvalues, which are always ± 1 or $\pm i$ as described below. In a real vector space, a unitary transformation can be thought of as simply a rigid rotation of the coordinate system, and all of the properties of a rigid rotation can be found in the unitary DFT.

The orthogonality of the DFT is now expressed as an [orthonormality](#) condition (which arises in many areas of mathematics as described in [root of unity](#)):

$$\sum_{m=0}^{N-1} U_{km} U_{mn}^* = \delta_{kn}$$

If \mathbf{X} is defined as the unitary DFT of the vector \mathbf{x} , then

$$X_k = \sum_{n=0}^{N-1} U_{kn} x_n$$

and the [Plancherel theorem](#) is expressed as

$$\sum_{n=0}^{N-1} x_n y_n^* = \sum_{k=0}^{N-1} X_k Y_k^*$$

If we view the DFT as just a coordinate transformation which simply specifies the components of a vector in a new coordinate system, then the above is just the statement that the dot product of two vectors is preserved under a unitary DFT transformation. For the special case $\mathbf{x} = \mathbf{y}$, this implies that the length of a vector is preserved as well—this is just [Parseval's theorem](#),

$$\sum_{n=0}^{N-1} |x_n|^2 = \sum_{k=0}^{N-1} |X_k|^2$$

A consequence of the [circular convolution theorem](#) is that the DFT matrix F diagonalizes any [circulant matrix](#).

Expressing the inverse DFT in terms of the DFT [\[edit\]](#)

A useful property of the DFT is that the inverse DFT can be easily expressed in terms of the (forward) DFT, via several well-known "tricks". (For example, in computations, it is often convenient to only implement a fast Fourier transform corresponding to one transform direction and then to get the other transform direction from the first.)

First, we can compute the inverse DFT by reversing the inputs (Duhamel *et al.*, 1988):

$$\mathcal{F}^{-1}(\{x_n\}) = \mathcal{F}(\{x_{N-n}\})/N$$

(As usual, the subscripts are interpreted [modulo](#) N ; thus, for $n = 0$, we have $x_{N-0} = x_0$.)

Second, one can also conjugate the inputs and outputs:

$$\mathcal{F}^{-1}(\mathbf{x}) = \mathcal{F}(\mathbf{x}^*)^*/N$$

Third, a variant of this conjugation trick, which is sometimes preferable because it requires no modification of the data values, involves swapping real and imaginary parts (which can be done on a computer simply by modifying [pointers](#)). Define $\text{swap}(x_n)$ as x_n with its real and imaginary parts swapped—that is, if $x_n = a + bi$ then $\text{swap}(x_n)$ is $b + ai$. Equivalently, $\text{swap}(x_n)$ equals $i x_n^*$. Then

$$\mathcal{F}^{-1}(\mathbf{x}) = \text{swap}(\mathcal{F}(\text{swap}(\mathbf{x}))) / N$$

That is, the inverse transform is the same as the forward transform with the real and imaginary parts swapped for both input and output, up to a normalization (Duhamel *et al.*, 1988).

The conjugation trick can also be used to define a new transform, closely related to the DFT, that is [involutory](#)—that is, which is its own inverse. In particular, $T(\mathbf{x}) = \mathcal{F}(\mathbf{x}^*)/\sqrt{N}$ is clearly its own inverse: $T(T(\mathbf{x})) = \mathbf{x}$. A closely related involutory transformation (by a factor of $(1+i)/\sqrt{2}$) is $H(\mathbf{x}) = \mathcal{F}((1+i)\mathbf{x}^*)/\sqrt{2N}$, since the $(1+i)$ factors in $H(H(\mathbf{x}))$ cancel the 2. For real inputs \mathbf{x} , the real part of $H(\mathbf{x})$ is none other than the [discrete Hartley transform](#), which is also involutory.

Eigenvalues and eigenvectors [edit]

The **eigenvalues** of the DFT matrix are simple and well-known, whereas the **eigenvectors** are complicated, not unique, and are the subject of ongoing research.

Consider the unitary form **U** defined above for the DFT of length *N*, where

$$U_{m,n} = \frac{1}{\sqrt{N}} \omega_N^{(m-1)(n-1)} = \frac{1}{\sqrt{N}} e^{-\frac{2\pi i}{N}(m-1)(n-1)}.$$

This matrix satisfies the **matrix polynomial** equation:

$$U^4 = I.$$

This can be seen from the inverse properties above: operating **U** twice gives the original data in reverse order, so operating **U** four times gives back the original data and is thus the **identity matrix**. This means that the eigenvalues λ satisfy the equation:

$$\lambda^4 = 1.$$

Therefore, the eigenvalues of **U** are the fourth **roots of unity**: λ is +1, −1, +i, or −i.

Since there are only four distinct eigenvalues for this $N \times N$ matrix, they have some **multiplicity**. The multiplicity gives the number of **linearly independent** eigenvectors corresponding to each eigenvalue. (Note that there are *N* independent eigenvectors; a unitary matrix is never **defective**.)

The problem of their multiplicity was solved by McClellan and Parks (1972), although it was later shown to have been equivalent to a problem solved by Gauss (Dickinson and Steiglitz, 1982). The multiplicity depends on the value of *N modulo 4*, and is given by the following table:

size <i>N</i>	$\lambda = +1$	$\lambda = -1$	$\lambda = -i$	$\lambda = +i$
$4m$	$m + 1$	m	m	$m - 1$
$4m + 1$	$m + 1$	m	m	m
$4m + 2$	$m + 1$	$m + 1$	m	m
$4m + 3$	$m + 1$	$m + 1$	$m + 1$	m

Multiplicities of the eigenvalues λ of the unitary DFT matrix **U as a function of the transform size *N* (in terms of an integer *m*).**

Otherwise stated, the **characteristic polynomial** of **U** is:

$$\det(\lambda I - U) = (\lambda - 1)^{\lfloor \frac{N+4}{4} \rfloor} (\lambda + 1)^{\lfloor \frac{N+2}{4} \rfloor} (\lambda + i)^{\lfloor \frac{N+1}{4} \rfloor} (\lambda - i)^{\lfloor \frac{N-1}{4} \rfloor}.$$

No simple analytical formula for general eigenvectors is known. Moreover, the eigenvectors are not unique because any linear combination of eigenvectors for the same eigenvalue is also an eigenvector for that eigenvalue. Various researchers have proposed different choices of eigenvectors, selected to satisfy useful properties like **orthogonality** and to have "simple" forms (e.g., McClellan and Parks, 1972; Dickinson and Steiglitz, 1982; Grünbaum, 1982; Atakishiyev and Wolf, 1997; Candan *et al.*, 2000; Hanna *et al.*, 2004; Gurevich and Hadani, 2008).

A straightforward approach is to discretize an eigenfunction of the continuous **Fourier transform**, of which the most famous is the **Gaussian function**. Since **periodic summation** of the function means discretizing its frequency spectrum and discretization means periodic summation of the spectrum, the discretized and periodically summed Gaussian function yields an eigenvector of the discrete transform:

$$\bullet F(m) = \sum_{k \in \mathbb{Z}} \exp\left(-\frac{\pi \cdot (m + N \cdot k)^2}{N}\right).$$

A closed form expression for the series is not known, but it converges rapidly.

Two other simple closed-form analytical eigenvectors for special DFT period *N* were found (Kong, 2008):

For DFT period $N = 2L + 1 = 4K + 1$, where *K* is an integer, the following is an eigenvector of DFT:

$$\bullet F(m) = \prod_{s=K+1}^L \left[\cos\left(\frac{2\pi}{N}m\right) - \cos\left(\frac{2\pi}{N}s\right) \right]$$

For DFT period $N = 2L = 4K$, where *K* is an integer, the following is an eigenvector of DFT:

$$\bullet F(m) = \sin\left(\frac{2\pi}{N}m\right) \prod_{s=K+1}^{L-1} \left[\cos\left(\frac{2\pi}{N}m\right) - \cos\left(\frac{2\pi}{N}s\right) \right]$$

The choice of eigenvectors of the DFT matrix has become important in recent years in order to define a discrete analogue of the **fractional Fourier transform**—the DFT matrix can be taken to fractional powers by exponentiating the eigenvalues (e.g., Rubio and Santhanam, 2005). For the **continuous Fourier transform**, the natural orthogonal eigenfunctions are the **Hermite functions**, so various discrete analogues of these have been employed as the eigenvectors of the DFT, such as the **Kravchuk polynomials** (Atakishiyev and Wolf, 1997). The "best" choice of eigenvectors to define a fractional discrete Fourier transform remains an open question, however.

Uncertainty principle [edit]

If the random variable X_k is constrained by

$$\sum_{n=0}^{N-1} |X_n|^2 = 1,$$

then

$$P_n = |X_n|^2$$

may be considered to represent a discrete **probability mass function** of *n*, with an associated probability mass function constructed from the transformed variable,

$$Q_m = N|x_m|^2.$$

For the case of continuous functions $P(x)$ and $Q(k)$, the **Heisenberg uncertainty principle** states that

$$D_0(X)D_0(x) \geq \frac{1}{16\pi^2}$$

where $D_0(X)$ and $D_0(x)$ are the variances of $|X|^2$ and $|x|^2$ respectively, with the equality attained in the case of a suitably normalized [Gaussian distribution](#). Although the variances may be analogously defined for the DFT, an analogous uncertainty principle is not useful, because the uncertainty will not be shift-invariant. Still, a meaningful uncertainty principle has been introduced by Massar and Spindel.^[6]

However, the Hirschman [entropic uncertainty](#) will have a useful analog for the case of the DFT.^[7] The Hirschman uncertainty principle is expressed in terms of the [Shannon entropy](#) of the two probability functions.

In the discrete case, the Shannon entropies are defined as

$$H(X) = - \sum_{n=0}^{N-1} P_n \ln P_n$$

and

$$H(x) = - \sum_{m=0}^{N-1} Q_m \ln Q_m ,$$

and the [entropic uncertainty](#) principle becomes^[7]

$$H(X) + H(x) \geq \ln(N) .$$

The equality is obtained for P_n equal to translations and modulations of a suitably normalized [Kronecker comb](#) of period A where A is any exact integer divisor of N . The probability mass function Q_m will then be proportional to a suitably translated [Kronecker comb](#) of period $B = N/A$.^[7]

There is also a well-known deterministic uncertainty principle that uses signal sparsity (or the number of non-zero coefficients).^[8] Let $\|x\|_0$ and $\|X\|_0$ be the number of non-zero elements of the time and frequency sequences x_0, x_1, \dots, x_{N-1} and X_0, X_1, \dots, X_{N-1} , respectively. Then,

$$N \leq \|x\|_0 \cdot \|X\|_0.$$

As an immediate consequence of the inequality of arithmetic and geometric means, one also has $2\sqrt{N} \leq \|x\|_0 + \|X\|_0$. Both uncertainty principles were shown to be tight for specifically-chosen "picket-fence" sequences (discrete impulse trains), and find practical use for signal recovery applications.^[8]

The real-input DFT ^[edit]

If x_0, \dots, x_{N-1} are [real numbers](#), as they often are in practical applications, then the DFT obeys the symmetry:

$$X_{N-k} \equiv X_{-k} = X_k^*, \text{ where } X^* \text{ denotes } \text{complex conjugation}.$$

It follows that X_0 and $X_{N/2}$ are real-valued, and the remainder of the DFT is completely specified by just $N/2-1$ complex numbers.

Generalized DFT (shifted and non-linear phase) ^[edit]

It is possible to shift the transform sampling in time and/or frequency domain by some real shifts a and b , respectively. This is sometimes known as a **generalized DFT** (or **GDFT**), also called the **shifted DFT** or **offset DFT**, and has analogous properties to the ordinary DFT:

$$X_k = \sum_{n=0}^{N-1} x_n e^{-\frac{2\pi i}{N}(k+b)(n+a)} \quad k = 0, \dots, N-1.$$

Most often, shifts of $1/2$ (half a sample) are used. While the ordinary DFT corresponds to a periodic signal in both time and frequency domains, $a = 1/2$ produces a signal that is anti-periodic in frequency domain ($X_{k+N} = -X_k$) and vice versa for $b = 1/2$. Thus, the specific case of $a = b = 1/2$ is known as an *odd-time odd-frequency* discrete Fourier transform (or O^2 DFT). Such shifted transforms are most often used for symmetric data, to represent different boundary symmetries, and for real-symmetric data they correspond to different forms of the discrete [cosine](#) and [sine](#) transforms.

Another interesting choice is $a = b = -(N-1)/2$ which is called the **centered DFT** (or **CDFT**). The centered DFT has the useful property that, when N is a multiple of four, all four of its eigenvalues (see above) have equal multiplicities (Rubio and Santhanam, 2005)^[9]

The term GDFT is also used for the non-linear phase extensions of DFT. Hence, GDFT method provides a generalization for constant amplitude orthogonal block transforms including linear and non-linear phase types. GDFT is a framework to improve time and frequency domain properties of the traditional DFT, e.g. auto/cross-correlations, by the addition of the properly designed phase shaping function (non-linear, in general) to the original linear phase functions (Akansu and Agirman-Tosun, 2010).^[10]

The discrete Fourier transform can be viewed as a special case of the [z-transform](#), evaluated on the unit circle in the complex plane; more general z-transforms correspond to *complex* shifts a and b above.

Multidimensional DFT ^[edit]

The ordinary DFT transforms a one-dimensional sequence or [array](#) x_n that is a function of exactly one discrete variable n . The multidimensional DFT of a multidimensional array x_{n_1, n_2, \dots, n_d} that is a function of d discrete variables $n_\ell = 0, 1, \dots, N_\ell - 1$ for ℓ in $1, 2, \dots, d$ is defined by:

$$X_{k_1, k_2, \dots, k_d} = \sum_{n_1=0}^{N_1-1} \left(\omega_{N_1}^{k_1 n_1} \sum_{n_2=0}^{N_2-1} \left(\omega_{N_2}^{k_2 n_2} \dots \sum_{n_d=0}^{N_d-1} \omega_{N_d}^{k_d n_d} \cdot x_{n_1, n_2, \dots, n_d} \right) \right) ,$$

where $\omega_{N_\ell} = \exp(-2\pi i/N_\ell)$ as above and the d output indices run from $k_\ell = 0, 1, \dots, N_\ell - 1$. This is more compactly expressed in [vector](#) notation, where we define $\mathbf{n} = (n_1, n_2, \dots, n_d)$ and $\mathbf{k} = (k_1, k_2, \dots, k_d)$ as d -dimensional vectors of indices from 0 to $\mathbf{N} - 1$, which we define as $\mathbf{N} - 1 = (N_1 - 1, N_2 - 1, \dots, N_d - 1)$:

$$X_{\mathbf{k}} = \sum_{\mathbf{n}=0}^{\mathbf{N}-1} e^{-2\pi i \mathbf{k} \cdot (\mathbf{n}/\mathbf{N})} x_{\mathbf{n}} ,$$

where the division \mathbf{n}/\mathbf{N} is defined as $\mathbf{n}/\mathbf{N} = (n_1/N_1, \dots, n_d/N_d)$ to be performed element-wise, and the sum denotes the set of nested

summations above.

The inverse of the multi-dimensional DFT is, analogous to the one-dimensional case, given by:

$$x_{\mathbf{n}} = \frac{1}{\prod_{\ell=1}^d N_{\ell}} \sum_{\mathbf{k}=0}^{\mathbf{N}-1} e^{2\pi i \mathbf{n} \cdot (\mathbf{k}/\mathbf{N})} X_{\mathbf{k}}.$$

As the one-dimensional DFT expresses the input $x_{\mathbf{n}}$ as a superposition of sinusoids, the multidimensional DFT expresses the input as a superposition of [plane waves](#), or multidimensional sinusoids. The direction of oscillation in space is \mathbf{k}/\mathbf{N} . The amplitudes are $X_{\mathbf{k}}$. This decomposition is of great importance for everything from [digital image processing](#) (two-dimensional) to solving [partial differential equations](#). The solution is broken up into plane waves.

The multidimensional DFT can be computed by the [composition](#) of a sequence of one-dimensional DFTs along each dimension. In the two-dimensional case x_{n_1, n_2} the N_1 independent DFTs of the rows (i.e., along n_2) are computed first to form a new array y_{n_1, k_2} . Then the N_2 independent DFTs of y along the columns (along n_1) are computed to form the final result X_{k_1, k_2} . Alternatively the columns can be computed first and then the rows. The order is immaterial because the nested summations above [commute](#).

An algorithm to compute a one-dimensional DFT is thus sufficient to efficiently compute a multidimensional DFT. This approach is known as the *row-column* algorithm. There are also intrinsically [multidimensional FFT algorithms](#).

The real-input multidimensional DFT [\[edit\]](#)

For input data x_{n_1, n_2, \dots, n_d} consisting of [real numbers](#), the DFT outputs have a conjugate symmetry similar to the one-dimensional case above:

$$X_{k_1, k_2, \dots, k_d} = X_{N_1 - k_1, N_2 - k_2, \dots, N_d - k_d}^*,$$

where the star again denotes complex conjugation and the ℓ -th subscript is again interpreted modulo N_{ℓ} (for $\ell = 1, 2, \dots, d$).

Applications [\[edit\]](#)

The DFT has seen wide usage across a large number of fields; we only sketch a few examples below (see also the references at the end). All applications of the DFT depend crucially on the availability of a fast algorithm to compute discrete Fourier transforms and their inverses, a [fast Fourier transform](#).

Spectral analysis [\[edit\]](#)

When the DFT is used for [signal spectral analysis](#), the $\{x_n\}$ sequence usually represents a finite set of uniformly spaced time-samples of some signal $x(t)$, where t represents time. The conversion from continuous time to samples (discrete-time) changes the underlying [Fourier transform](#) of $x(t)$ into a [discrete-time Fourier transform](#) (DTFT), which generally entails a type of distortion called [aliasing](#). Choice of an appropriate sample-rate (see [Nyquist rate](#)) is the key to minimizing that distortion. Similarly, the conversion from a very long (or infinite) sequence to a manageable size entails a type of distortion called [leakage](#), which is manifested as a loss of detail (a.k.a. resolution) in the DTFT. Choice of an appropriate sub-sequence length is the primary key to minimizing that effect. When the available data (and time to process it) is more than the amount needed to attain the desired frequency resolution, a standard technique is to perform multiple DFTs, for example to create a [spectrogram](#). If the desired result is a power spectrum and noise or randomness is present in the data, averaging the magnitude components of the multiple DFTs is a useful procedure to reduce the [variance](#) of the spectrum (also called a [periodogram](#) in this context); two examples of such techniques are the [Welch method](#) and the [Bartlett method](#); the general subject of estimating the power spectrum of a noisy signal is called [spectral estimation](#).

A final source of distortion (or perhaps *illusion*) is the DFT itself, because it is just a discrete sampling of the DTFT, which is a function of a continuous frequency domain. That can be mitigated by increasing the resolution of the DFT. That procedure is illustrated at [Sampling the DTFT](#).

- The procedure is sometimes referred to as *zero-padding*, which is a particular implementation used in conjunction with the [fast Fourier transform](#) (FFT) algorithm. The inefficiency of performing multiplications and additions with zero-valued "samples" is more than offset by the inherent efficiency of the FFT.
- As already noted, leakage imposes a limit on the inherent resolution of the DTFT. So there is a practical limit to the benefit that can be obtained from a fine-grained DFT.

Filter bank [\[edit\]](#)

See [FFT filter banks](#) and [Sampling the DTFT](#).

Data compression [\[edit\]](#)

The field of digital signal processing relies heavily on operations in the frequency domain (i.e. on the Fourier transform). For example, several [lossy](#) image and sound compression methods employ the discrete Fourier transform: the signal is cut into short segments, each is transformed, and then the Fourier coefficients of high frequencies, which are assumed to be unnoticeable, are discarded. The decompressor computes the inverse transform based on this reduced number of Fourier coefficients. (Compression applications often use a specialized form of the DFT, the [discrete cosine transform](#) or sometimes the [modified discrete cosine transform](#).) Some relatively recent compression algorithms, however, use [wavelet transforms](#), which give a more uniform compromise between time and frequency domain than obtained by chopping data into segments and transforming each segment. In the case of [JPEG2000](#), this avoids the spurious image features that appear when images are highly compressed with the original [JPEG](#).

Partial differential equations [\[edit\]](#)

Discrete Fourier transforms are often used to solve [partial differential equations](#), where again the DFT is used as an approximation for the [Fourier series](#) (which is recovered in the limit of infinite N). The advantage of this approach is that it expands the signal in complex exponentials e^{inx} , which are eigenfunctions of differentiation: $d/dx e^{inx} = in e^{inx}$. Thus, in the Fourier representation, differentiation is simple—we just multiply by in . (Note, however, that the choice of n is not unique due to aliasing; for the method to be convergent, a choice similar to that in the [trigonometric interpolation](#) section above should be used.) A [linear differential equation](#) with constant coefficients is transformed into an easily solvable algebraic equation. One then uses the inverse DFT to transform the result back into the ordinary spatial representation. Such an approach is called a [spectral method](#).

Polynomial multiplication [\[edit\]](#)

Suppose we wish to compute the polynomial product $c(x) = a(x) \cdot b(x)$. The ordinary product expression for the coefficients of c involves a linear (acyclic) convolution, where indices do not "wrap around." This can be rewritten as a cyclic convolution by taking the coefficient vectors for $a(x)$ and

$b(x)$ with constant term first, then appending zeros so that the resultant coefficient vectors **a** and **b** have dimension $d > \deg(a(x)) + \deg(b(x))$. Then,

c = **a** * **b**

Where **c** is the vector of coefficients for $c(x)$, and the convolution operator $*$ is defined so

$$c_n = \sum_{m=0}^{d-1} a_m b_{n-m \bmod d} \qquad n = 0, 1 \dots, d-1$$

But convolution becomes multiplication under the DFT:

$$\mathcal{F}(\mathbf{c}) = \mathcal{F}(\mathbf{a})\mathcal{F}(\mathbf{b})$$

Here the vector product is taken elementwise. Thus the coefficients of the product polynomial $c(x)$ are just the terms $0, \dots, \deg(a(x)) + \deg(b(x))$ of the coefficient vector

$$\mathbf{c} = \mathcal{F}^{-1}(\mathcal{F}(\mathbf{a})\mathcal{F}(\mathbf{b})).$$

With a [fast Fourier transform](#), the resulting algorithm takes $O(N \log N)$ arithmetic operations. Due to its simplicity and speed, the [Cooley–Tukey FFT algorithm](#), which is limited to [composite](#) sizes, is often chosen for the transform operation. In this case, d should be chosen as the smallest integer greater than the sum of the input polynomial degrees that is factorizable into small prime factors (e.g. 2, 3, and 5, depending upon the FFT implementation).

Multiplication of large integers [\[edit\]](#)

The fastest known [algorithms](#) for the multiplication of very large [integers](#) use the polynomial multiplication method outlined above. Integers can be treated as the value of a polynomial evaluated specifically at the number base, with the coefficients of the polynomial corresponding to the digits in that base. After polynomial multiplication, a relatively low-complexity carry-propagation step completes the multiplication.

Convolution [\[edit\]](#)

When data is [convolved](#) with a function with wide support, such as for downsampling by a large sampling ratio, because of the [Convolution theorem](#) and the FFT algorithm, it may be faster to transform it, multiply pointwise by the transform of the filter and then reverse transform it. Alternatively, a good filter is obtained by simply truncating the transformed data and re-transforming the shortened data set.

Some discrete Fourier transform pairs [\[edit\]](#)

Some DFT pairs		
$x_n = \frac{1}{N} \sum_{k=0}^{N-1} X_k e^{i2\pi kn/N}$	$X_k = \sum_{n=0}^{N-1} x_n e^{-i2\pi kn/N}$	Note
$x_n e^{i2\pi n\ell/N}$	$X_{k-\ell}$	Shift theorem
$x_{n-\ell}$	$X_k e^{-i2\pi k\ell/N}$	
$x_n \in \mathbb{R}$	$X_k = X_{N-k}^*$	Real DFT
a^n	$\begin{cases} N & \text{if } a = e^{i2\pi k/N} \\ \frac{1-a^N}{1-a e^{-i2\pi k/N}} & \text{otherwise} \end{cases}$	from the geometric progression formula
$\binom{N-1}{n}$	$(1 + e^{-i2\pi k/N})^{N-1}$	from the binomial theorem
$\begin{cases} \frac{1}{W} & \text{if } 2n < W \text{ or } 2(N-n) < W \\ 0 & \text{otherwise} \end{cases}$	$\begin{cases} 1 & \text{if } k = 0 \\ \frac{\sin(\frac{\pi W k}{N})}{W \sin(\frac{\pi k}{N})} & \text{otherwise} \end{cases}$	x_n is a rectangular window function of W points centered on $n=0$, where W is an odd integer , and X_k is a sinc-like function (specifically, X_k is a Dirichlet kernel)
$\sum_{j \in \mathbb{Z}} \exp\left(-\frac{\pi}{cN} \cdot (n + N \cdot j)^2\right)$	$\sqrt{cN} \cdot \sum_{j \in \mathbb{Z}} \exp\left(-\frac{\pi c}{N} \cdot (k + N \cdot j)^2\right)$	Discretization and periodic summation of the scaled Gaussian functions for $c > 0$. Since either c or $\frac{1}{c}$ is larger than one and thus warrants fast convergence of one of the two series, for large c you may choose to compute the frequency spectrum and convert to the time domain using the discrete Fourier transform.

Generalizations [\[edit\]](#)

Representation theory [\[edit\]](#)

For more details on this topic, see [Representation theory of finite groups § Discrete Fourier transform](#).

The DFT can be interpreted as the complex-valued [representation theory](#) of the finite [cyclic group](#). In other words, a sequence of n complex numbers can be thought of as an element of n -dimensional complex space \mathbf{C}^n or equivalently a function f from the finite cyclic group of order n to the complex numbers, $\mathbf{Z}_n \rightarrow \mathbf{C}$. So f is a [class function](#) on the finite cyclic group, and thus can be expressed as a linear combination of the irreducible characters of this group, which are the roots of unity.

From this point of view, one may generalize the DFT to representation theory generally, or more narrowly to the [representation theory of finite groups](#).

More narrowly still, one may generalize the DFT by either changing the target (taking values in a field other than the complex numbers), or the domain (a group other than a finite cyclic group), as detailed in the sequel.

Other fields [\[edit\]](#)

Main articles: [Discrete Fourier transform \(general\)](#) and [Number-theoretic transform](#)

Many of the properties of the DFT only depend on the fact that $e^{-\frac{2\pi i}{N}}$ is a [primitive root of unity](#), sometimes denoted ω_N or W_N (so that $\omega_N^N = 1$

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External links [edit]

- Interactive explanation of the DFT
- Matlab tutorial on the Discrete Fourier Transformation
- Interactive flash tutorial on the DFT
- Mathematics of the Discrete Fourier Transform by Julius O. Smith III
- Fast implementation of the DFT - coded in C and under General Public License (GPL)
- The DFT "à Pied": Mastering The Fourier Transform in One Day
- Explained: The Discrete Fourier Transform
- wavetable Cooker GPL application with graphical interface written in C, and implementing DFT IDFT to generate a wavetable set
- Online DFT coefficients calculator with graphical display – Gives online calculation of DFT coefficients with graphical display
- Discrete Fourier Transform JavaScript Program – A live, online, program for computing the discrete Fourier Transform coefficients of a real, uniformly-spaced, data sequence.

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Sub-fields	Audio signal processing · Digital image processing · Speech processing · Statistical signal processing	
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