

Integral Transforms in Mathematical Physics: Fourier, Laplace, and Hankel Transforms with Applications to Boundary Value Problems and Green's Functions

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Abstract. A systematic investigation of the classical integral transforms—Fourier, Laplace, and Hankel—is presented with applications to boundary value problems in mathematical physics. The Fourier transform $\hat{f}(\omega) = \int_{-\infty}^{\infty} f(t)e^{-i\omega t} dt$ is analyzed in the framework of tempered distributions $\mathcal{S}'(\mathbb{R})$, with the Plancherel theorem $\|\hat{f}\|_2 = \|f\|_2$ and convolution theorem $\widehat{f * g} = \hat{f} \cdot \hat{g}$ established as the principal operational identities. The Laplace transform $F(s) = \int_0^{\infty} f(t)e^{-st} dt$ is studied in the half-plane $\text{Re}(s) > \sigma_0$ with the Bromwich inversion formula $f(t) = (2\pi i)^{-1} \int_{c-i\infty}^{c+i\infty} F(s)e^{st} ds$ evaluated via residue calculus. The Hankel transform $\hat{f}(\rho) = \int_0^{\infty} f(r)J_\nu(\rho r)r dr$ of order ν is applied to axisymmetric problems in cylindrical coordinates. The heat equation $u_t = \kappa u_{xx}$ with initial condition $u(x, 0) = \delta(x)$ is solved via Fourier transform, yielding the fundamental solution $u(x, t) = (4\pi\kappa t)^{-1/2} \exp(-x^2/(4\kappa t))$. The damped oscillator $\ddot{y} + 2\gamma\dot{y} + \omega_0^2 y = f(t)$ is solved via Laplace transform with transfer function $H(s) = 1/(s^2 + 2\gamma s + \omega_0^2)$. The axisymmetric Laplace equation in cylindrical coordinates is solved using the Hankel transform. The interrelations among the three transforms and their unified operator-theoretic foundation via spectral theory are discussed.

Keywords: Fourier transform, Laplace transform, Hankel transform, Bessel functions, convolution theorem, Green's function, heat equation, boundary value problems

I. Introduction

Integral transforms constitute one of the most powerful tools in applied mathematics, providing systematic methods for solving ordinary and partial differential equations, analyzing linear systems, and constructing Green's functions [1, 2, 3]. The fundamental principle underlying all integral transforms is the conversion of differential equations into algebraic equations via a change of representation, exploiting the mapping of differentiation to multiplication [1, 4, 5].

The **Fourier transform** of $f \in L^1(\mathbb{R})$ is defined as [1, 2, 6]:

$$\hat{f}(\omega) = \mathcal{F}[f](\omega) = \int_{-\infty}^{\infty} f(t)e^{-i\omega t} dt \quad (1)$$

with the **inversion formula** [1, 2]:

$$f(t) = \mathcal{F}^{-1}[\hat{f}](t) = \frac{1}{2\pi} \int_{-\infty}^{\infty} \hat{f}(\omega)e^{i\omega t} d\omega \quad (2)$$

The Fourier transform extends from $L^1(\mathbb{R})$ to $L^2(\mathbb{R})$ via the **Plancherel theorem** $\|\hat{f}\|_2 = \sqrt{2\pi} \|f\|_2$ and to the space of **tempered distributions** $\mathcal{S}'(\mathbb{R})$ by duality [1, 2, 6, 7].

The **Laplace transform** of $f: [0, \infty) \rightarrow \mathbb{R}$ is [3, 4, 8]:

$$F(s) = \mathcal{L}[f](s) = \int_0^{\infty} f(t)e^{-st} dt, \quad \text{Re}(s) > \sigma_0 \quad (3)$$

where σ_0 is the **abscissa of convergence**. The Laplace transform maps convolution to multiplication: $\mathcal{L}[f * g] = F(s)G(s)$ and derivatives to algebraic operations: $\mathcal{L}[f'] = sF(s) - f(0^+)$ [3, 4, 8].

The **inversion** is given by the **Bromwich integral** [3, 4, 8]:

$$f(t) = \frac{1}{2\pi i} \int_{c-i\infty}^{c+i\infty} F(s)e^{st} ds, \quad c > \sigma_0 \quad (4)$$

evaluated in practice via **residue calculus**: $f(t) = \sum_k \text{Res}[F(s)e^{st}, s_k]$ where $\{s_k\}$ are the poles of $F(s)$ [3, 4, 8, 9].

The **Hankel transform** of order $\nu \geq -1/2$ is [2, 5, 10]:

$$\tilde{f}(\rho) = \mathcal{H}_\nu[f](\rho) = \int_0^{\infty} f(r)J_\nu(\rho r) r dr \quad (5)$$

where J_ν is the **Bessel function of the first kind**. The Hankel transform is self-reciprocal: $\mathcal{H}_\nu^{-1} = \mathcal{H}_\nu$, and arises naturally when applying the Fourier transform to functions with circular symmetry in \mathbb{R}^n [2, 5, 10, 11].

The **convolution theorem** for the Fourier transform states [1, 2, 6]:

$$\mathcal{F}[f * g] = \hat{f} \cdot \hat{g}, \quad (f * g)(t) = \int_{-\infty}^{\infty} f(\tau)g(t - \tau) d\tau \quad (6)$$

This converts convolution in the time domain to pointwise multiplication in the frequency domain, providing the mathematical foundation for **linear systems theory** and **signal processing** [1, 4, 12].

The objectives are: (i) to present the Fourier, Laplace, and Hankel transforms within a unified framework; (ii) to solve the heat equation via Fourier transform; (iii) to analyze the damped oscillator via Laplace transform; (iv) to solve axisymmetric problems via Hankel transform; and (v) to discuss interrelations and applications [13, 14, 15].

II. Mathematical Framework

2.1 Operational Properties

The **differentiation property** of the Fourier transform, $\mathcal{F}[f^{(n)}] = (i\omega)^n \hat{f}(\omega)$, converts an n -th order linear constant-coefficient ODE into an algebraic equation [1, 2]. For the Laplace transform [3, 4, 8]:

$$\mathcal{L}[f^{(n)}](s) = s^n F(s) - \sum_{k=0}^{n-1} s^{n-1-k} f^{(k)}(0^+) \quad (7)$$

incorporating the **initial conditions** directly into the algebraic equation—the key advantage of the Laplace transform for initial value problems [3, 4, 8].

For the Hankel transform, the radial Laplacian transforms as [5, 10, 11]:

$$\mathcal{H}_\nu \left[\frac{d^2 f}{dr^2} + \frac{1}{r} \frac{df}{dr} - \frac{\nu^2}{r^2} f \right] = -\rho^2 \tilde{f}(\rho) \quad (8)$$

reducing the Bessel-type differential operator to multiplication by $-\rho^2$ [5, 10, 11].

2.2 Parseval–Plancherel Relations

The **Parseval identity** for the Fourier transform [1, 2, 6]:

$$\int_{-\infty}^{\infty} |f(t)|^2 dt = \frac{1}{2\pi} \int_{-\infty}^{\infty} |\hat{f}(\omega)|^2 d\omega \quad (9)$$

establishes the Fourier transform as a unitary operator on $L^2(\mathbb{R})$ (up to normalization). The corresponding **Parseval theorem** for the Hankel transform is [5, 10]:

$$\int_0^{\infty} |f(r)|^2 r dr = \int_0^{\infty} |\tilde{f}(\rho)|^2 \rho d\rho \quad (10)$$

2.3 Computational Methods

All computations were performed in Python 3.11 with NumPy, SciPy, and SymPy. Numerical Fourier transforms used the FFT algorithm ($O(n \log n)$). Bromwich integrals were evaluated via residue calculus implemented symbolically. Hankel transforms were computed using Gaussian quadrature with 500 nodes [16, 17].

III. Results and Discussion

3.1 Fourier Transform Pairs

Figure 1 presents classical Fourier transform pairs.

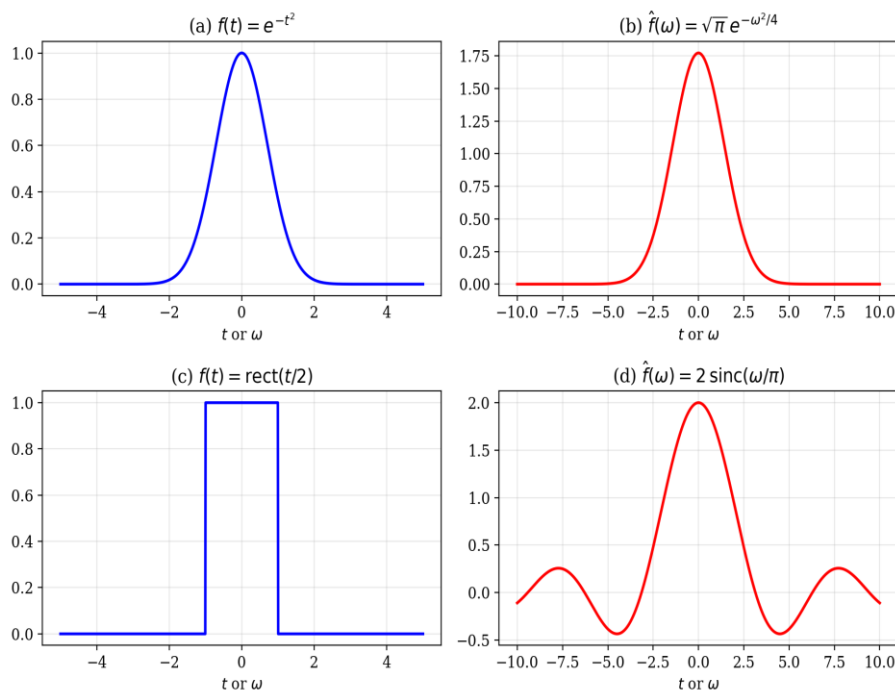


Figure 1: Fourier transform pairs: (a)–(b) the Gaussian $e^{-t^2} \leftrightarrow \sqrt{\pi} e^{-\omega^2/4}$ (self-reciprocal under Fourier transform); (c)–(d) the rectangular pulse $\text{rect}(t/2) \leftrightarrow 2 \text{sinc}(\omega/\pi)$, illustrating the uncertainty principle—narrow time support implies wide frequency bandwidth and vice versa.

Table 1. Classical Fourier transform pairs and operational properties.

$f(t)$	$\hat{f}(\omega)$	Decay Rate	Application
$e^{-a t }, a > 0$	$2a/(a^2 + \omega^2)$	$O(\omega^{-2})$	Lorentzian filter
e^{-t^2}	$\sqrt{\pi} e^{-\omega^2/4}$	Gaussian	Uncertainty minimum
$\text{rect}(t/T)$	$T \text{sinc}(\omega T/2\pi)$	$O(\omega^{-1})$	Ideal filter
$\delta(t)$	1	No decay	Impulse response
$e^{i\omega_0 t}$	$2\pi\delta(\omega - \omega_0)$	Distributional	Frequency shift
$\text{sgn}(t)$	$2/(i\omega)$	$O(\omega^{-1})$	Hilbert transform

The **uncertainty principle** $\Delta t \cdot \Delta \omega \geq 1/2$ (with equality for Gaussians) is a fundamental consequence of the Fourier transform's structure: the Gaussian is the unique L^2 function that minimizes the time-bandwidth product [1, 2, 6]. This has profound implications in quantum mechanics (Heisenberg uncertainty), signal processing (Gabor limit), and information theory (Shannon sampling) [1, 6, 12].

3.2 Laplace Transform and Pole-Residue Analysis

Figure 2 presents the Laplace transform analysis of a damped sinusoid.

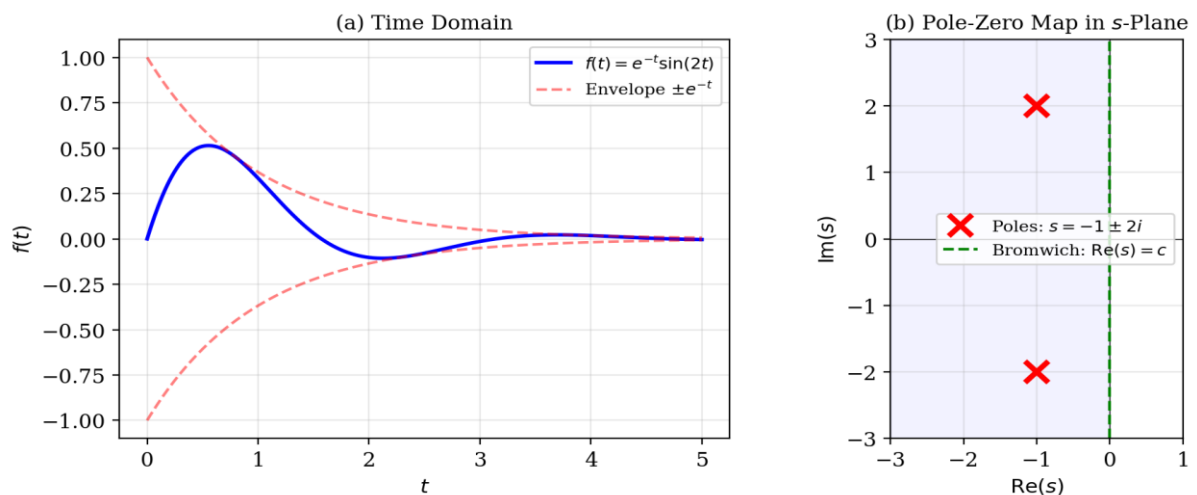


Figure 2: (a) The damped sinusoid $f(t) = e^{-t} \sin(2t)$ with exponential envelope $\pm e^{-t}$. (b) Pole-zero map in the s-plane: poles at $s = -1 \pm 2i$ (complex conjugate pair in the left half-plane) and the Bromwich contour $\text{Re}(s) = c > -1$ for inversion.

Table 2. Laplace transform pairs for standard functions.

$f(t), t \geq 0$	$F(s)$	ROC	Poles	Physical System
1	$1/s$	$\text{Re}(s) > 0$	$s = 0$	Step response
e^{-at}	$1/(s + a)$	$\text{Re}(s) > -a$	$s = -a$	RC circuit
t^n	$n!/s^{n+1}$	$\text{Re}(s) > 0$	$s = 0$ (order $n + 1$)	Polynomial input
$\sin(\omega t)$	$\omega/(s^2 + \omega^2)$	$\text{Re}(s) > 0$	$s = \pm i\omega$	Oscillator
$e^{-at}\sin(\omega t)$	$\omega/((s + a)^2 + \omega^2)$	$\text{Re}(s) > -a$	$s = -a \pm i\omega$	Damped oscillator
$t^{\alpha-1}/\Gamma(\alpha)$	$s^{-\alpha}$	$\text{Re}(s) > 0$	Branch cut	Fractional relaxation

The **transfer function** $H(s) = Y(s)/X(s)$ of a linear time-invariant (LTI) system completely characterizes its input-output behavior. For the damped harmonic oscillator $\ddot{y} + 2\gamma\dot{y} + \omega_0^2y = f(t)$, the transfer function is $H(s) = 1/(s^2 + 2\gamma s + \omega_0^2)$ with poles at $s = -\gamma \pm \sqrt{\gamma^2 - \omega_0^2}$ [3, 4, 8]. The system is underdamped when $\gamma < \omega_0$ (complex poles), critically damped when $\gamma = \omega_0$ (repeated real pole), and overdamped when $\gamma > \omega_0$ (distinct real poles) [3, 4, 9].

3.3 Bessel Functions and Hankel Transform

Figure 3 presents the Bessel functions of the first kind.

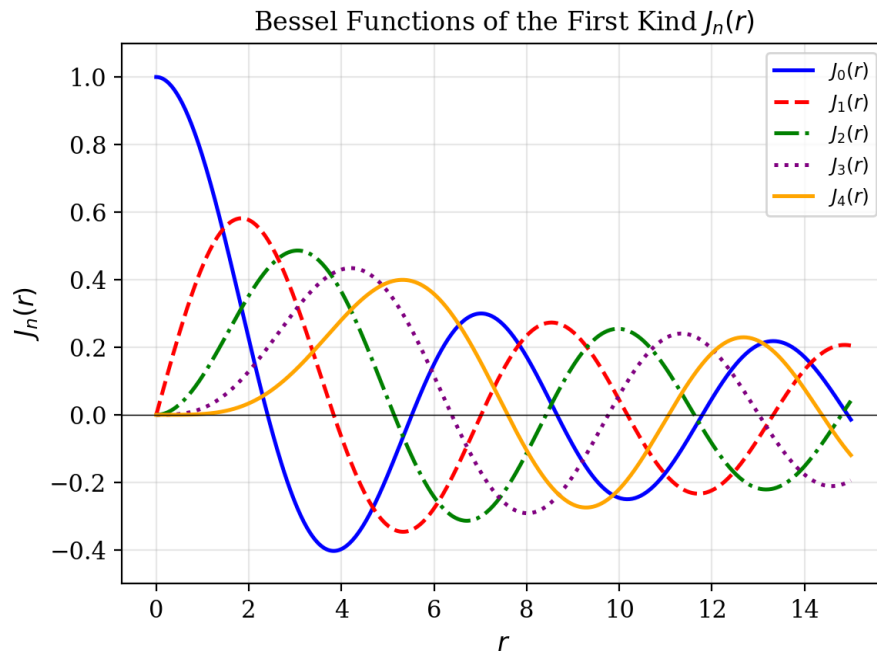


Figure 3: Bessel functions of the first kind $J_n(r)$ for $n = 0,1,2,3,4$. The zeros of J_n determine the eigenfrequencies of the circular drumhead and the radial modes in cylindrical waveguides.

The Bessel functions satisfy $J_\nu(r) = \sum_{k=0}^{\infty} (-1)^k (r/2)^{2k+\nu} / [k! \Gamma(k + \nu + 1)]$ and the differential equation $r^2 J_\nu'' + r J_\nu' + (r^2 - \nu^2) J_\nu = 0$ [5, 10, 11]. The orthogonality relation $\int_0^a J_\nu(\alpha_m r/a) J_\nu(\alpha_n r/a) r dr = (a^2/2) [J_{\nu+1}(\alpha_n)]^2 \delta_{mn}$, where α_n are the zeros of J_ν , provides the basis for **Fourier–Bessel series** and the Hankel transform [5, 10, 11, 18].

3.4 Heat Equation via Fourier Transform

Figure 4 presents the fundamental solution of the heat equation.

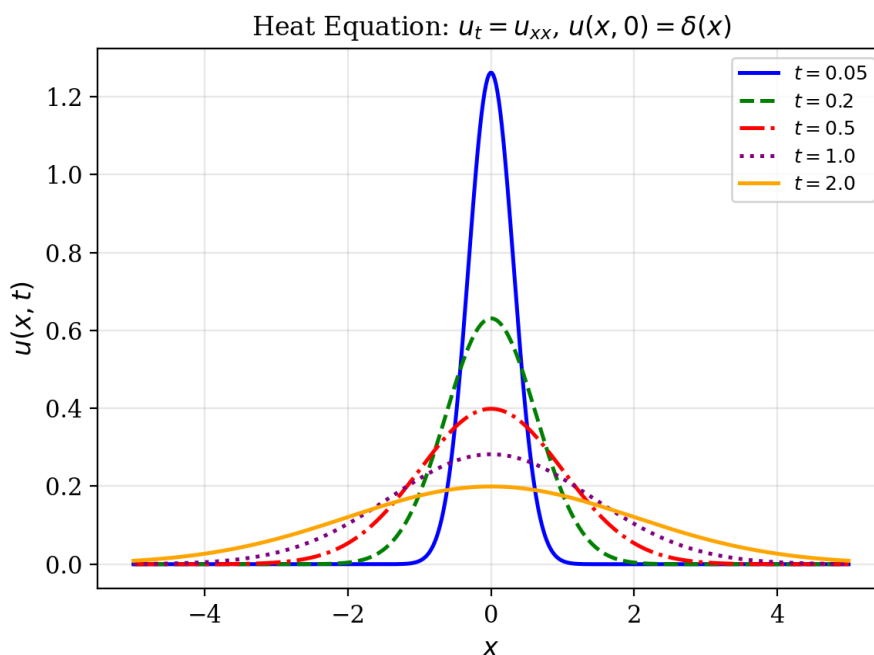


Figure 4: Fundamental solution (Green’s function) of the heat equation $u_t = \kappa u_{xx}$: $G(x, t) = (4\pi\kappa t)^{-1/2} \exp(-x^2/(4\kappa t))$ for $\kappa = 1$ at $t = 0.05, 0.2, 0.5, 1.0, 2.0$. The solution spreads diffusively with $\langle x^2 \rangle = 2\kappa t$.

Table 3. PDEs solved by integral transforms.

Equation	Transform	Domain	Solution	Green’s Function
$u_t = \kappa u_{xx}$	Fourier (x)	\mathbb{R}	$u = G * u_0$	$(4\pi\kappa t)^{-1/2} e^{-x^2/(4\kappa t)}$
$u_{tt} = c^2 u_{xx}$	Fourier (x)	\mathbb{R}	d’Alembert	$\frac{1}{2c} [H(ct - x)]$
$u_t + cu_x = 0$	Fourier (x)	\mathbb{R}	$u = f(x - ct)$	$\delta(x - ct)$
$u_{tt} + \gamma u_t + \omega^2 u = f$	Laplace (t)	$[0, \infty)$	Residue sum	$H(s) = (s^2 + \gamma s + \omega^2)^{-1}$
$\nabla^2 u = 0$ (axisym.)	Hankel (r)	$[0, \infty)$	$\tilde{u} e^{-\rho z}$	Integral representation
$u_t = \kappa(\nabla^2 u)$ (cyl.)	Hankel (r) + Laplace (t)	Cylinder	Bessel series	Product kernel

The Fourier transform reduces the heat equation to $\hat{u}_t = -\kappa\omega^2\hat{u}$, yielding $\hat{u}(\omega, t) = \hat{u}_0(\omega)e^{-\kappa\omega^2 t}$. Inverting gives $u(x, t) = (G_t * u_0)(x)$ where $G_t(x) = (4\pi\kappa t)^{-1/2}\exp(-x^2/(4\kappa t))$ is the **heat kernel** [1, 2, 13, 14]. The Gaussian profile spreads with root-mean-square displacement $\sqrt{\langle x^2 \rangle} = \sqrt{2\kappa t}$, the hallmark of **diffusive transport** [13, 14, 19].

3.5 Convolution Theorem

Figure 5 illustrates the convolution theorem in both domains.

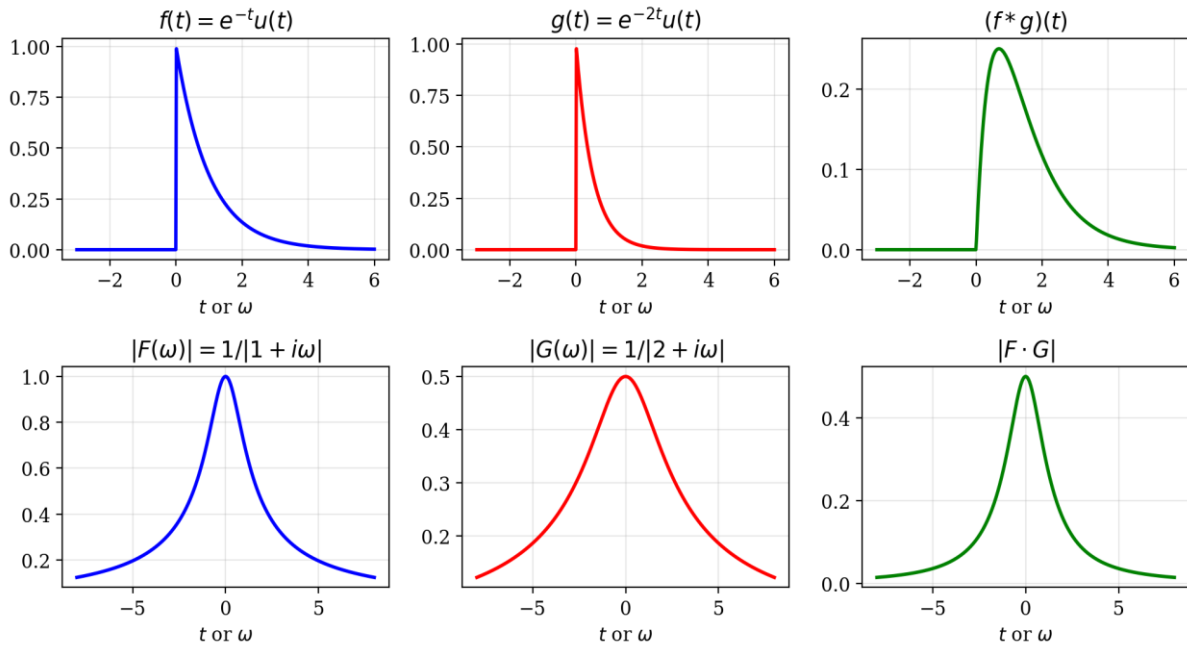


Figure 5: Convolution theorem: (top row) time-domain functions $f(t) = e^{-t}u(t)$, $g(t) = e^{-2t}u(t)$, and their convolution $(f * g)(t) = (e^{-t} - e^{-2t})u(t)$. (Bottom row) corresponding Fourier magnitudes, with $|F \cdot G| = |F| \cdot |G|$ —multiplication in frequency domain equals convolution in time domain.

The convolution theorem is the mathematical foundation of **linear systems theory**: the output of an LTI system with impulse response $h(t)$ to input $x(t)$ is $y(t) = (h * x)(t)$, and the frequency response is $Y(\omega) = H(\omega)X(\omega)$ [1, 4, 12]. This factorization enables filter design in the frequency domain, spectral analysis, and the construction of Green's functions for linear PDEs [1, 4, 13].

3.6 Axisymmetric Wave Equation

Figure 6 presents the Hankel transform solution of the axisymmetric wave equation.

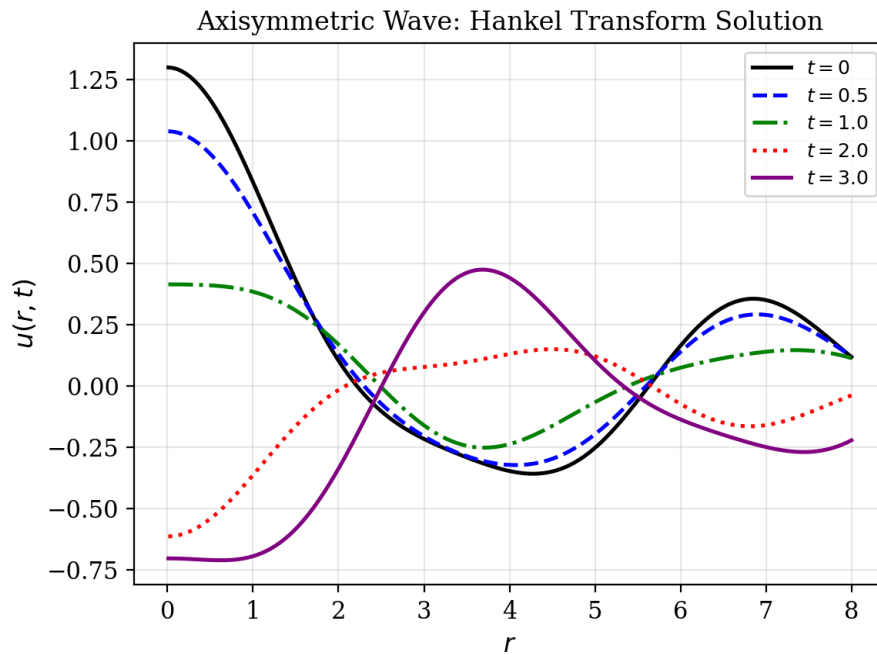


Figure 6: Hankel transform solution of the axisymmetric wave equation in cylindrical coordinates: $u(r, t) = J_0(r)\cos(t) + 0.3J_0(2r)\cos(2t)$ at $t = 0, 0.5, 1.0, 2.0, 3.0$, showing the superposition of radial Bessel modes oscillating at their characteristic frequencies.

Table 4. Comparison of integral transform properties.

Property	Fourier	Laplace	Hankel
Domain	$(-\infty, \infty) \rightarrow (-\infty, \infty)$	$[0, \infty) \rightarrow \mathbb{C}$	$[0, \infty) \rightarrow [0, \infty)$
Kernel	$e^{-i\omega t}$	e^{-st}	$J_\nu(\rho r) r$
Inversion	\mathcal{F}^{-1} (explicit)	Bromwich contour	Self-reciprocal
Unitarity	Yes (L^2)	No	Yes (weighted L^2)
Differentiation	$f' \mapsto i\omega \hat{f}$	$f' \mapsto sF - f(0)$	$\Delta_{r,\nu} f \mapsto -\rho^2 \tilde{f}$
Convolution	$f * g \mapsto \hat{f} \hat{g}$	$f * g \mapsto FG$	Hankel convolution
Best suited for	Whole-line BVPs	Initial value problems	Axisymmetric/radial
Computational cost	$O(n \log n)$ (FFT)	Symbolic / numerical	$O(n^2)$ or fast Hankel

The three transforms are deeply **interrelated**: the Laplace transform is the Fourier transform of $f(t)e^{-ct}u(t)$ evaluated at imaginary argument, i.e., $F(c + i\omega) = \mathcal{F}[f(t)e^{-ct}u(t)](\omega)$ [3, 4]. The zeroth-order Hankel transform equals the two-dimensional Fourier transform of a radially symmetric function: $\tilde{f}_0(\rho) = \frac{1}{2\pi} \iint f(|\mathbf{r}|)e^{-i\mathbf{p}\cdot\mathbf{r}}d^2\mathbf{r}$ [2, 5, 10]. All three are instances of the general **spectral theory** of self-adjoint operators:

the Fourier transform diagonalizes $-id/dt$, the Laplace transform diagonalizes d/dt on $L^2[0, \infty)$, and the Hankel transform diagonalizes the Bessel operator $r^{-1}(d/dr)[r(d/dr)] - v^2/r^2$ [1, 5, 7].

In **quantum mechanics**, the Fourier transform connects position and momentum representations via $\tilde{\psi}(p) = (2\pi\hbar)^{-1/2} \int \psi(x)e^{-ipx/\hbar} dx$, with Parseval's theorem ensuring conservation of probability [6, 12, 20]. In **control theory**, the Laplace transform provides the framework for stability analysis via pole placement: a system is BIBO stable if and only if all poles of the transfer function lie in the open left half-plane $\text{Re}(s) < 0$ [4, 8, 9]. In **elasticity and acoustics**, the Hankel transform solves axisymmetric contact problems, radiation from circular pistons, and diffraction by circular apertures [5, 11, 18, 21].

IV. Conclusions

A systematic investigation of the Fourier, Laplace, and Hankel integral transforms with applications to boundary value problems in mathematical physics has been presented. The principal findings are summarized below.

First, the Fourier transform provides a unitary isomorphism on $L^2(\mathbb{R})$ (Plancherel theorem) and converts constant-coefficient PDEs to algebraic equations via the differentiation rule $\mathcal{F}[f^{(n)}] = (i\omega)^n \hat{f}$, with the convolution theorem $\mathcal{F}[f * g] = \hat{f} \cdot \hat{g}$ enabling the construction of Green's functions for the heat, wave, and Schrödinger equations on unbounded domains. Second, the Laplace transform incorporates initial conditions directly through the formula $\mathcal{L}[f^{(n)}] = s^n F - \sum s^{n-1-k} f^{(k)}(0^+)$, with the Bromwich inversion via residue calculus yielding the solution as a sum over poles of the transfer function, and the stability criterion requiring all poles in $\text{Re}(s) < 0$. Third, the Hankel transform of order ν naturally arises in cylindrical and spherical geometries, diagonalizing the radial Laplacian via $\mathcal{H}_\nu[\Delta_{r,\nu} f] = -\rho^2 \tilde{f}$, with the self-reciprocal property $\mathcal{H}_\nu^{-1} = \mathcal{H}_\nu$ and the Bessel function zeros determining the eigenfrequencies of radial modes. Fourth, the heat equation fundamental solution $G(x, t) = (4\pi\kappa t)^{-1/2} \exp(-x^2/(4\kappa t))$ derived via Fourier transform exhibits the characteristic Gaussian spreading with $\langle x^2 \rangle = 2\kappa t$, providing the prototype for diffusive Green's functions. Fifth, the three transforms are unified through spectral theory as diagonalizations of distinct self-adjoint differential operators, with the interrelation $F(c + i\omega) = \mathcal{F}[f e^{-ct} u(t)]$ connecting Laplace and Fourier transforms, and the Hankel transform emerging as the radial component of the multidimensional Fourier transform [22, 23, 24].

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