"Normal modes" in Physics equations

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Maxwell's Equations:

$$\nabla \cdot \mathbf{E} = \frac{\rho}{\varepsilon_0}$$

$$\nabla \cdot \mathbf{B} = 0$$

$$\nabla \times \mathbf{E} = -\frac{\partial \mathbf{B}}{\partial t}$$

$$\nabla \times \mathbf{B} = \mu_0 \mathbf{J} + \mu_0 \varepsilon_0 \frac{\partial \mathbf{E}}{\partial t}$$
(1)
(2)
(3)

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Maxwell's Equations

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$$\nabla \times \mathbf{E} = -\frac{\partial \mathbf{B}}{\partial t} \tag{3}$$

$$\nabla \times \mathbf{B} = \mu_0 \mathbf{J} + \mu_0 \varepsilon_0 \frac{\partial \mathbf{E}}{\partial t}$$
 (4)

Consider no charges $\rho=0$ and no currents J=0. With $c=\frac{1}{\sqrt{\mu_0\varepsilon_0}}$ this gives the vector wave equations:

$$abla^2 \mathbf{E} - \frac{1}{c^2} \, \partial_t^2 \mathbf{E} = \mathbf{0}, \qquad
abla^2 \mathbf{B} - \frac{1}{c^2} \, \partial_t^2 \mathbf{B} = \mathbf{0}.$$

Consider a parallel-plate cavity on $x \in [0, L]$ and a transverse field component $E_y(x, t) = \varphi(t) \phi(x)$.

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The 1D wave equation becomes

$$\phi''(x) + k^2 \phi(x) = 0,$$
 $\ddot{\varphi}(t) + c^2 k^2 \varphi(t) = 0.$

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At a perfect electric conductor the tangential electric field vanishes: For plates at x = 0, L this gives

$$E_y(0,t) = E_y(L,t) = 0 \Rightarrow \phi(0) = \phi(L) = 0$$
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Eigenmodes. (Normal modes) $\phi_n(x) = \sin(\frac{n\pi x}{L})$, $\varphi_n(t) = \exp(\pm i\omega_n t)$, $k_n = \frac{n\pi}{L}$, $\omega_n = ck_n$, $n = 1, 2, \ldots$

Finite Differences

On the computer the derivative can be approximated as:

$$\frac{\partial \phi}{\partial x}(x_i) \approx \frac{\phi(x_{i+1}) - \phi(x_{i-1})}{2\Delta x}$$

$$\frac{\partial^2 \phi}{\partial x^2}(x_i) \approx \frac{\phi(x_{i+1}) - 2\phi(x_i) + \phi(x_{i-1})}{\Delta x^2}$$
(5)

$$\frac{\partial^2 \phi}{\partial x^2}(x_i) \approx \frac{\phi(x_{i+1}) - 2\phi(x_i) + \phi(x_{i-1})}{\Delta x^2} \tag{6}$$

where we assume to have a grid discretization $x_0, x_1, \dots, x_N, x_{N+1}, \Delta x = x_i - x_{i-1}$ and the function evaluated on those points $\phi(x_i) = \phi_i$.

Finite Differences

Then considering the derivative approximation + BC we obtain:

$$\frac{1}{\Delta x^{2}} \begin{bmatrix}
-2 & 1 & 0 & \cdots & 0 & 0 \\
1 & -2 & 1 & \cdots & 0 & 0 \\
0 & 1 & -2 & \cdots & 0 & 0 \\
\vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\
0 & 0 & 0 & \cdots & 1 & -2
\end{bmatrix} \begin{bmatrix}
\phi_{1} \\
\phi_{2} \\
\phi_{3} \\
\vdots \\
\phi_{N-1} \\
\phi_{N}
\end{bmatrix} = -k^{2} \begin{bmatrix}
\phi_{1} \\
\phi_{2} \\
\phi_{3} \\
\vdots \\
\phi_{N-1} \\
\phi_{N}
\end{bmatrix} (7)$$

where we use the boundary condition $\phi_0 = 0$ and $\phi_{N+1} = 0$.

Wave Equation Eigensystem Code

```
function analytical_eigenstate(i_mode)
    return x -> sin(x * pi * i_mode/ p.L)
end
T = Float64
p = (L = 1,
   N = 100.
    c = 1
A = make_EM_wave_Dirichlet_matrix(T, p)
x = LinRange(T(0), T(p.L), p.N+2)
F = eigen(Matrix(A), sortby=abs)
```

Wave Equation solution

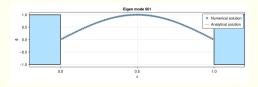


Figure 1: First eigen mode

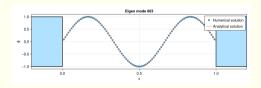


Figure 2: Third eigen mode

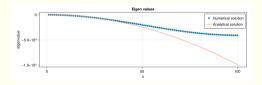


Figure 3: Wave Equation Eigenvalues

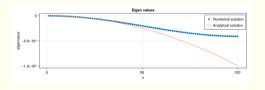


Figure 3: Wave Equation Eigenvalues

Can we do better?

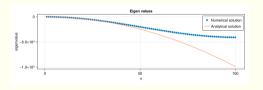


Figure 3: Wave Equation Eigenvalues

Can we do better?

• Higher order finite difference approx

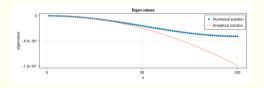


Figure 3: Wave Equation Eigenvalues

Can we do better?

- Higher order finite difference approx
- Finite Element Method

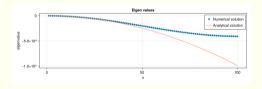


Figure 3: Wave Equation Eigenvalues

Can we do better?

- Higher order finite difference approx
- Finite Element Method
- Sine Fourier Transform

Diffusion Equation

Diffusion Equation

Diffusion Equation:

$$\frac{\partial \rho}{\partial t} = \nabla \cdot (D\nabla \rho) \tag{8}$$

- ullet ho particle density.
- $\mathbf{j} = -D\nabla \rho$ particle density current.

Consider a parallel-walls cavity on $x \in [0, L]$, a probability density $\rho(x, t) = \varphi(t) \phi(x)$, and D constant.

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The 1D diffusion equation becomes

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The 1D diffusion equation becomes

$$\phi''(x) + k^2 \phi(x) = 0, \qquad \dot{\varphi}(t) + Dk^2 \varphi(t) = 0.$$

Eigenmodes. (Normal modes)
$$\phi_n(x) = \cos(\frac{n\pi x}{L})$$
, $\varphi_n(t) = \exp(-\tau_n t)$, $k_n = \frac{n\pi}{L}$, $\tau_n = Dk_n^2$, $n = 0, 1, \ldots$

Finite Differences

On the computer the derivative + Neumann BC can be approximated as:

$$\frac{1}{\Delta x^{2}} \begin{bmatrix}
-1 & 1 & 0 & \cdots & 0 & 0 \\
1 & -2 & 1 & \cdots & 0 & 0 \\
0 & 1 & -2 & \cdots & 0 & 0 \\
\vdots & \vdots & \vdots & \ddots & \vdots & \vdots \\
0 & 0 & 0 & \cdots & 1 & -1
\end{bmatrix} \begin{bmatrix}
\phi_{1} \\
\phi_{2} \\
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\end{bmatrix} = -k^{2} \begin{bmatrix}
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\phi_{2} \\
\phi_{3} \\
\vdots \\
\phi_{N-1} \\
\phi_{N}
\end{bmatrix} (9)$$

where we use the boundary condition $\nabla \phi_0 = 0$ and $\nabla \phi_{N+1} = 0$.

Difussion Equation Eigensystem Code

```
function analytical_eigenstate(i_mode)
    return x -> cos(x * pi * i_mode/ p.L)
end
T = Float64
p = (L = 1,
   N = 100.
    c = 1
A = make_Diffusion_eq_Neumann_matrix(T, p)
x = LinRange(T(0), T(p.L), p.N+2)
F = eigen(Matrix(A), sortby=abs)
```

Diffusion Equation solution

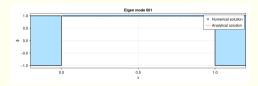


Figure 4: First eigen mode

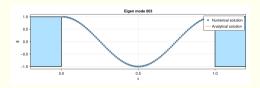


Figure 5: Third eigen mode

Schrödinger Equation in position basis:

$$i\hbar \frac{\partial \psi}{\partial t} = -\frac{\hbar^2}{2m} \nabla^2 \psi + V(\mathbf{x}) \psi \tag{10}$$

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- ullet ψ gives probability amplitude.
- $-\frac{\hbar^2}{2m}\nabla^2\psi$ comes from momentum contribution to the Hamiltonian.
- V(x) is the potential energy contribution to the Hamiltonian.

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- ullet ψ gives probability amplitude.
- $-\frac{\hbar^2}{2m}\nabla^2\psi$ comes from momentum contribution to the Hamiltonian.
- V(x) is the potential energy contribution to the Hamiltonian.

We will consider $V(\mathbf{x}) = V_0$ for $x \in [-L, L]$ and $V(\mathbf{x}) = 0$ elsewhere.

Finite well $V_0 < 0$

Bound states

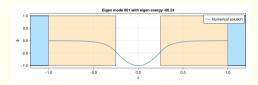


Figure 6: First bound state

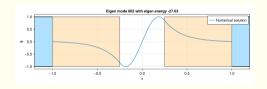


Figure 7: Second bound state

Finite well $V_0 < 0$

No bound states

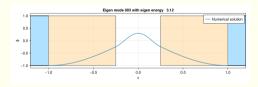


Figure 8: First no bound state

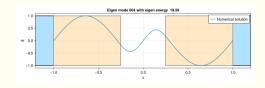


Figure 9: Second no bound state

Finite barrier $V_0 > 0$

Near degenerated eigen states

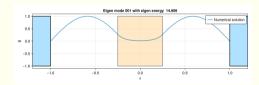


Figure 10: First eigen state

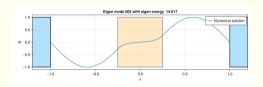


Figure 11: Second eigen state

Sum of near degenerated eigen states