

DD2370 Computational Methods for Electromagnetics

Eigenvalue Problems in CEM

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Eigenvalues Problems

- CEM applications can be either
 1. a driven system, where one seeks the **response to a source**, for instance an antenna, or
 2. an eigenvalue calculation, where one seeks the **natural oscillation frequencies of the system**.
- The field solution to the eigenvalue calculation is a nonzero field that satisfies the homogeneous problem, i.e. the field problem without sources.
 - This type of situation occurs typically for an electromagnetic system with a source that is nonzero for a finite time.
 - Once the source vanishes, the field solution can be expressed as a superposition of eigenmodes, where each eigenmode oscillates at a particular eigenfrequency.

Maxwell's Equations in Frequency Domain

In a linear, dispersion-free medium, Maxwell's equations can be written as the single second-order curl-curl equation for the electric field

$$-\epsilon \frac{\partial^2 \mathbf{E}}{\partial t^2} = \nabla \times \frac{1}{\mu} \nabla \times \mathbf{E} + \frac{\partial \mathbf{J}}{\partial t} \quad \longrightarrow \quad \boxed{\omega_m^2 \epsilon \mathbf{E}_m = \nabla \times \frac{1}{\mu} \nabla \times \mathbf{E}_m}$$

In the absence of sources, $\mathbf{J}=0$, and with harmonic time dependence $\exp(jt)$, the curl-curl equation gives the eigenvalue problem

Eigenvalues and Eigenfunctions

- For nontrivial solutions ($\mathbf{E}_m \neq 0$)
 - ω_m = eigenvalue
 - \mathbf{E}_m is the corresponding eigenfunction, or eigenmode.
- If the region is a closed cavity with a perfectly conducting boundary (i.e., $\mathbf{n} \times \mathbf{E} = 0$ on the boundaries), the operator on the right-hand side, $\nabla \times \frac{1}{\mu} \nabla \times \mathbf{E}_m$, is **self-adjoint**, that is

$$\omega_m^2 \epsilon \mathbf{E}_m = \nabla \times \frac{1}{\mu} \nabla \times \mathbf{E}_m$$

$$\int_{\Omega} \mathbf{E}_1 \cdot L[\mathbf{E}_2] dV = \int_{\Omega} \mathbf{E}_2 \cdot L[\mathbf{E}_1] dV$$

Eigenvalues in Closed Cavity with PEC

- We can calculate analytically the eigenvalues as

$$\omega^2 = \frac{\int_{\Omega} \mu^{-1} |\nabla \times \mathbf{E}|^2 dV}{\int_{\Omega} \epsilon |\mathbf{E}|^2 dV},$$

- Eigenvalues are **real** and **nonnegative**.
- The eigenfrequencies are real for any lossless region bounded by perfect conductors.
 - Damping can appear if there is dissipation of energy, for example from regions with finite electrical conductivity, or if the region is not completely enclosed by a perfect conductor.

Helmholtz Equation

The self-adjoint curl-curl equation leads us to consider eigenvalue problems of the type

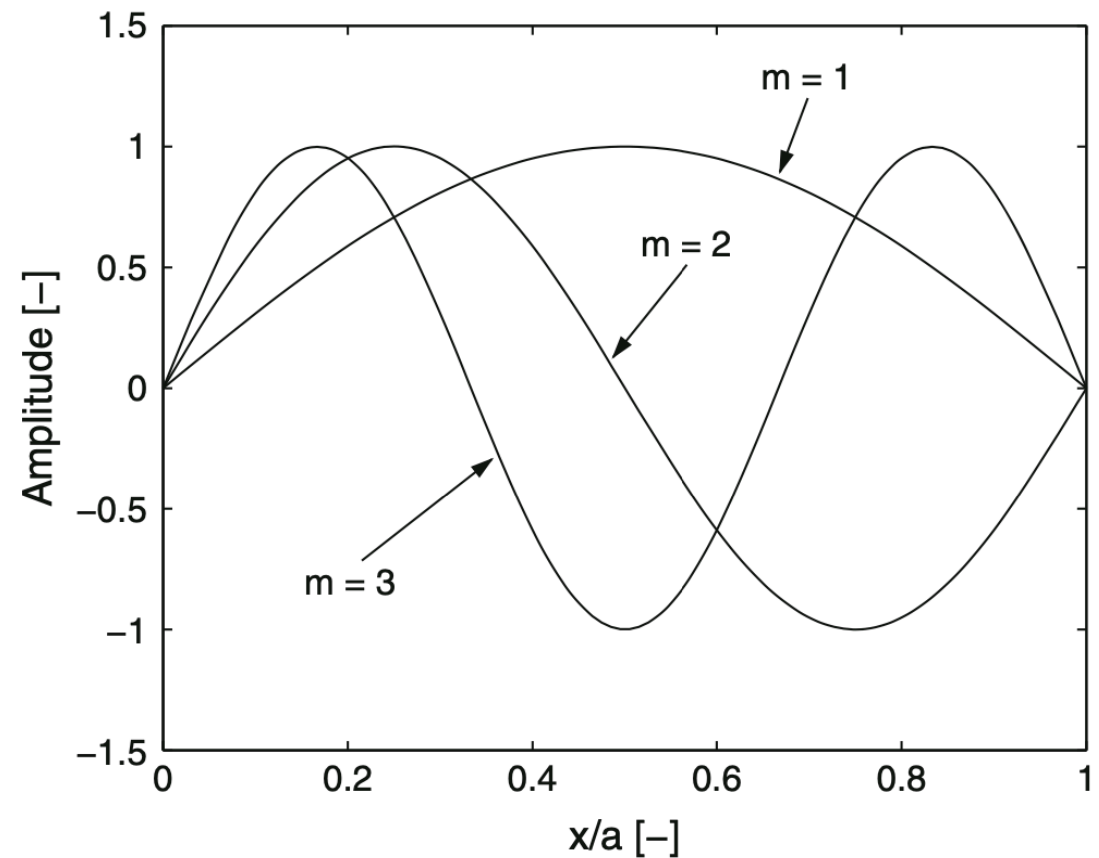
$$L[f] = -\omega^2 f \quad \text{in } \Omega$$

together with a suitable boundary condition on $\partial\Omega$ and a . We will assume that L is a linear self-adjoint operator with nonpositive eigenvalues. As a simple example to illustrate general principles, we will study the 1D **Helmholtz equation**:

$$\frac{d^2 f}{dx^2} = -k^2 f, \quad 0 < x < a, \quad f(0) = f(a) = 0$$

Solutions of Eigenvalue Problem

$$k_m = \frac{m\pi}{a}, \quad m \text{ an integer,}$$



Solve Numerically Helmholtz Equation with FD

To discretize the 1D Helmholtz equation by finite differences, we divide the interval $[0, a]$ into N subintervals of equal length $h = a/N$. The simplest finite difference approximation is

$$\frac{d^2 f}{dx^2} = -k^2 f, \quad 0 < x < a, \quad f(0) = f(a) = 0$$



$$\frac{f_{i-1} - 2f_i + f_{i+1}}{h^2} = -k^2 f_i, \quad i = 1, 2, \dots, N-1$$

Matrix Formulation of Eigenvalue Problem

The boundary conditions are $f_0 = f_N = 0$, so there is no reason to include f_0 and f_N as unknowns.

$\frac{f_{i-1} - 2f_i + f_{i+1}}{h^2} = -k^2 f_i$ can be written as a linear system with an $(N - 1) \times (N - 1)$ matrix **A**:

$$\mathbf{A}\mathbf{f} = \lambda\mathbf{f}$$

$$\mathbf{A} = \frac{1}{h^2} \begin{pmatrix} -2 & 1 & 0 & 0 & 0 & 0 \\ 1 & -2 & 1 & 0 & 0 & 0 \\ 0 & 1 & -2 & 1 & 0 & 0 \\ 0 & 0 & 1 & -2 & 1 & 0 \\ 0 & 0 & 0 & 1 & -2 & 1 \\ 0 & 0 & 0 & 0 & 1 & -2 \end{pmatrix}$$

Matlab Solution – Sparse Representation and *eig*

- When n is large, **A** consists mostly of zeros, and this can be exploited by saving the matrix in sparse form.
 - We can use **spalloc** function
- The physical eigenvalues $-k^2$ are simply the eigenvalues of the matrix **A**.
 - These eigenvalues can be computed with the MATLAB routine **eig**, which computes all eigenvalues and corresponding eigenvectors of an algebraic eigenvalue problem.

Matlab Code

```
% -----  
% Compute eigenvalues of 1D Helmholtz equation using FD  
% -----  
function k = HFD1D(a, N)  
  
% Arguments:  
%   a = length of interval  
%   N = number of subintervals (equal length)  
% Returns:  
%   k = eigenvalues  
  
h = a/N; % Grid size  
A = spalloc(N-1, ... % Allocate sparse matrix  
            N-1, ... % with 3*(N-1) nonzeros  
            3*(N-1));  
  
d = -2/h^2; % Value of diagonal entries  
s = 1/h^2; % Value of upper and lower  
            % diagonal entries  
  
% Initialize the diagonal entries  
for i = 1:N-1  
    A(i,i) = d; % Diagonal entries  
end  
  
% Initialize the upper and lower diagonal entries  
for i = 1:N-2  
    A(i,i+1) = s; % Upper diagonal entries  
    A(i+1,i) = s; % Lower diagonal entries  
end  
  
% Computing the eigenvalues  
lambda = eig(A);  
k = sqrt(sort(-lambda));
```

Calculate eigenvalues in $[0, \pi]$

```
>> k = HFD1D(pi,10)
```

```
k =
```

```
0.9959  
1.9673  
2.8902  
3.7420  
4.5016  
5.1504  
5.6723  
6.0546  
6.2878
```

```
>> k = HFD1D(pi,20)
```

```
k =
```

```
0.9990  
1.9918  
2.9723  
3.9345  
4.8725  
5.7804  
6.6527  
7.4839  
8.2690  
...
```

```
>> k = HFD1D(pi,30)
```

```
k =
```

```
0.9995  
1.9963  
2.9877  
3.9708  
4.9431  
5.9018  
6.8443  
7.7681  
8.6706  
9.5493  
...
```

We can perform convergence tests and find second order convergence