

The Laplacian Matrix and Spectral Graph Drawing. Courant-Fischer.



Zhiping (Patricia) Xiao
University of California, Los Angeles

October 8, 2020

Introduction

Resources

Basic Problems Examples

Background

Eigenvalues and Optimization: The Courant-Fischer Theorem

The Laplacian and Graph Drawing

Introduction



Course: Spectral Graph Theory from Yale.

Textbooks include:

- ▶ Spectral and Algebraic Graph Theory (Daniel A. Spielman)
- ▶ Scalable Algorithms for Data and Network Analysis (Shang-Hua Teng)

Objective of the course:

- ▶ To explore what eigenvalues and eigenvectors of graphs can tell us about their structure.

Prerequisites:

- ▶ Linear algebra, graphs, etc.

Textbook chapters:

- ▶ Spectral and Algebraic Graph Theory (Daniel A. Spielman) Chap 1 ~ 3
- ▶ Scalable Algorithms for Data and Network Analysis (Shang-Hua Teng) Chap 2.4

Problems listed in Prof. Teng's book Chap 2.4

- ▶ Significant Nodes: Ranking and Centrality
- ▶ Coherent Groups: Clustering and Communities
- ▶ Interplay between Networks and Dynamic Processes
- ▶ Multiple Networks: Composition and Similarity

Identifying nodes of relevance and significance. e.g.:

*Which nodes are the most **significant** nodes in a network or a sub-network? How quickly can we identify them?*

Significance could be measured either *numerically*, or by *ranking* the nodes.

Network centrality is a form of “dimensionality reduction” from “high dimensional” network data to “low dimensional” centrality measures or rankings.

e.g. PageRank

Identifying groups with significant structural properties.

Fundamental questions include:

- ▶ What are the significant clusters in a data set?
- ▶ How fast can we identify one, uniformly sample one, or enumerate all significant groups?
- ▶ How should we evaluate the consistency of a clustering or community-identification scheme?
- ▶ What desirable properties should clustering or community identification schemes satisfy?

Understanding the interplay between dynamic processes and their underlying networks.

A given social network can be part of different dynamic processes (e.g. epidemic spreading, viral marketing), which can potentially affect the relations between nodes. Fundamental questions include:

- ▶ How should we model the interaction between network nodes in a given dynamic process?
- ▶ How should we characterize node significance and group coherence with respect to a dynamic process?
- ▶ How fast can we identify influential nodes and significant communities?

To understand multiple networks instead of individual networks.

- ▶ network composition, e.g. multi-layer social network, multi-view graphs
- ▶ network similarity
 - ▶ similarity between two different networks
 - ▶ construct a sparser network that approximates a known one

$G = (V, E)$ (Friendship graphs, Network graphs, Circuit graphs, Protein-Protein Interaction graphs, etc.)

- ▶ G : a graph/network
- ▶ V : its vertex/node set
- ▶ E : its edge set (pair of vertices); edges have weight 1 by default, could assign other weights optionally.

By default (unless otherwise specified), a graph to be discussed will be:

- ▶ undirected (unordered vertices pairs in E)
- ▶ simple (having no loops or multiple edges)
- ▶ finite (V and E being finite sets)

Why we care about matrices?

Given a vector $\mathbf{x} \in \mathbb{R}^n$ and a matrix $\mathbf{M} \in \mathbb{R}^{n \times m}$

- ▶ \mathbf{M} could be an operator: $\mathbf{M}\mathbf{x} \in \mathbb{R}^m$
- ▶ \mathbf{M} could be used to define a quadratic form: $\mathbf{x}^T \mathbf{M} \mathbf{x} \in \mathbb{R}$

Adjacency matrix \mathbf{M}_G of $G = (V, E)$:

$$\mathbf{M}_G(a, b) = \begin{cases} 1 & \text{if } (a, b) \in E \\ 0 & \text{otherwise} \end{cases}$$

- ▶ most natural matrix to associate with a graph
- ▶ least useful

*This statement is made because it is only a **spread-sheet**, neither a natural **operator** or a natural **quadratic form**.*

Diffusion operator \mathbf{D}_G of $G = (V, E)$ is a diagonal matrix, probably the most natural operator associated with G :

$$\mathbf{D}_G(a, a) = \mathbf{d}(a)$$

where $\mathbf{d}(a)$ is the degree of vertex a .

- ▶ unweighted case: number of edges attached to it
- ▶ weighted case: weighted degree

$$\mathbf{d} \stackrel{\text{def}}{=} \mathbf{M}_G \mathbf{1}$$

There is a linear operator \mathbf{W}_G defined as:

$$\mathbf{W}_G = \mathbf{M}_G \mathbf{D}_G^{-1}$$

regarded as an operator denoting the *changes* of the graph between time steps.

Recall that diffusion operator \mathbf{D}_G is a diagonal matrix, \mathbf{W}_G is merely a rescaling of \mathbf{M}_G if the graph is *regular*¹.

With vector $\mathbf{p} \in \mathbb{R}^n$ denoting the values of n vertices (called “*distribution of how much stuff*” in the textbook), the distribution of stuff at each vertex will be $\mathbf{W}_G \mathbf{p}$.

¹Regular graph's vertices have the same degree.

$$\mathbf{W}_G = \mathbf{M}_G \mathbf{D}_G^{-1}$$

The next time step is:

$$\mathbf{W}_G \mathbf{p} = \mathbf{M}_G \mathbf{D}_G^{-1} \mathbf{p}$$

Think about the case where \mathbf{p} is a one-hot vector δ_a where only $\delta_a(a) = 1$ and all other elements are 0.

$$\mathbf{W}_G \delta_a = \mathbf{M}_G \mathbf{D}_G^{-1} \delta_a = \mathbf{M}_G (\mathbf{D}_G^{-1} \delta_a)$$

We find the vector $\mathbf{D}_G^{-1} \delta_a$ has value $1/\mathbf{d}(a)$ at vertex a and 0 everywhere else; $\mathbf{M}_G \mathbf{D}_G^{-1} \delta_a$ has value $1/\mathbf{d}(a)$ at all a 's **neighbors** and 0 otherwise.

A commonly-seen form of \mathbf{W}_G is sometimes more convenient:

$$\widetilde{\mathbf{W}}_G = I/2 + \mathbf{W}_G/2$$

describing a *lazy random walk* (1/2 chance stay, 1/2 chance go).

One of the purposes of spectral theory is to understand what happens when a linear operator like \mathbf{W}_G is repeatedly applied.

Laplacian matrix \mathbf{L}_G , the most natural quadratic form associated with the graph G :

$$\mathbf{L}_G \stackrel{\text{def}}{=} \mathbf{D}_G - \mathbf{M}_G$$

Given a vector $\mathbf{x} \in \mathbb{R}^n$, who could also be viewed as a *function* over the vertices, we have: ²

$$\mathbf{x}^T \mathbf{L}_G \mathbf{x} = \sum_{(a,b) \in E} w_{a,b} (\mathbf{x}(a) - \mathbf{x}(b))^2$$

representing the Laplacian quadratic form of a weighted graph ($w_{a,b}$ is the weight of edge (a,b)), could be used to measures the smoothness of \mathbf{x} (it would be small if \mathbf{x} is not changing drastically over any edge).

²Note that G has to be undirected

An example ($w_{a,b} = 1$):

$$\mathbf{M}_G = \begin{bmatrix} 0 & 1 & 1 \\ 1 & 0 & 0 \\ 1 & 0 & 0 \end{bmatrix} \quad \mathbf{D}_G = \begin{bmatrix} 2 & 0 & 0 \\ 0 & 1 & 0 \\ 0 & 0 & 1 \end{bmatrix}$$

$$\mathbf{L}_G = \mathbf{D}_G - \mathbf{M}_G = \begin{bmatrix} 2 & -1 & -1 \\ -1 & 1 & 0 \\ -1 & 0 & 1 \end{bmatrix}$$

$$\begin{aligned} \mathbf{x}^T \mathbf{L}_G \mathbf{x} &= x_1(2x_1 - x_2 - x_3) + x_2(-x_1 + x_2) + x_3(-x_1 + x_3) \\ &= 2x_1^2 + x_2^2 + x_3^2 - 2x_1x_2 - 2x_1x_3 = (x_1 - x_2)^2 + (x_1 - x_3)^2 \end{aligned}$$

*Intuitively, \mathbf{L}_G , \mathbf{D}_G and \mathbf{M}_G could be viewed as the sum of many subgraphs, each containing **one** edge.*

Review: the spectral theory for symmetric matrices (or those similar to symmetric matrices).

\mathbf{A} is similar to \mathbf{B} if there exists non-singular \mathbf{X} such that $\mathbf{X}^{-1}\mathbf{A}\mathbf{X} = \mathbf{B}$.

A vector ψ is an eigenvector of a matrix \mathbf{M} with eigenvalue λ if:

$$\mathbf{M}\psi = \lambda\psi$$

λ is an eigenvalue if and only if $\lambda\mathbf{I} - \mathbf{M}$ is a singular matrix ($\therefore \det(\lambda\mathbf{I} - \mathbf{M}) = 0$). The eigenvalues are the **roots** of the *characteristic polynomial* of \mathbf{M} :

$$\det(x\mathbf{I} - \mathbf{M})$$

in other words, being a solution to the *characteristic equation*:

$$\det(x\mathbf{I} - \mathbf{M}) = 0$$

Additional explanation on why “ λ is an eigenvalue if and only if $\lambda\mathbf{I} - \mathbf{M}$ is a singular matrix”: ³

$$\mathbf{M}\psi = \lambda\psi$$

$$(\lambda\mathbf{I} - \mathbf{M})\psi = 0$$

is a homogeneous linear system for ψ , with a trivial zero solution ($\psi = 0$).

A homogeneous linear system has a nonzero solution $\psi \neq 0$ iff its coefficient matrix (in this case, $\lambda\mathbf{I} - \mathbf{M}$), is singular.

³https://www-users.math.umn.edu/~olver/num_/lnv.pdf

Theorem (1.3.1 The Spectral Theorem)

*If \mathbf{M} is an n -by- n , real, symmetric matrix, **then** there exist real numbers $\lambda_1, \dots, \lambda_n$ and n mutually orthogonal unit vectors ψ_1, \dots, ψ_n and such that ψ_i is an eigenvector of \mathbf{M} of eigenvalue λ_i , for each i .*

If the matrix \mathbf{M} is not symmetric, it might not have n eigenvalues. And, even if it has n eigenvalues, their eigenvectors will not be orthogonal (linearly independent). Many studies will no longer apply to it when the matrix is not symmetric.

Review: solving the eigenvalues and eigenvectors. ⁴

$$\mathbf{M} = \begin{bmatrix} 0 & 1 \\ -2 & -3 \end{bmatrix}$$

$$\mathbf{M}\psi = \lambda\psi$$

$$(\mathbf{M} - \lambda\mathbf{I})\psi = 0$$

The determinant value of $\mathbf{M} - \lambda\mathbf{I}$ is 0 (by definition of the singular matrix, etc.).

$$\det(\mathbf{M} - \lambda\mathbf{I}) = 0$$

$$\det \left(\begin{bmatrix} -\lambda & 1 \\ -2 & -3 - \lambda \end{bmatrix} \right) = \lambda^2 + 3\lambda + 2 = (\lambda + 1)(\lambda + 2) = 0$$

The eigenvalues are:

$$\lambda_1 = -1, \lambda_2 = -2$$

Next we want to find the corresponding eigenvectors ψ_1 and ψ_2 , by solving:

$$(\mathbf{M} - \lambda\mathbf{I})\psi = 0$$

which means,

$$\begin{bmatrix} -\lambda_i & 1 \\ -2 & -3 - \lambda_i \end{bmatrix} \begin{bmatrix} \psi_{i,1} \\ \psi_{i,2} \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \end{bmatrix}$$

$$\psi_{i,2} - \lambda_i \psi_{i,1} = 0$$

$$2\psi_{i,1} + (3 + \lambda_i)\psi_{i,2} = 0$$

With $\lambda_1 = -1$, we have:

$$\begin{aligned}\psi_{1,2} + \psi_{1,1} &= 0 \\ 2\psi_{1,1} + 2\psi_{1,2} &= 0\end{aligned}$$

so the only constraint is that $\psi_{1,2} = -\psi_{1,1}$. We can choose any arbitrary constant k_1 and make it:

$$\psi_1 = k_1 \begin{bmatrix} 1 \\ -1 \end{bmatrix}$$

With $\lambda_2 = -2$, we have:

$$\psi_{2,2} + 2\psi_{2,1} = 0$$

$$2\psi_{2,1} + \psi_{2,2} = 0$$

again, we need an arbitrary constant k_2 and we have:

$$\psi_2 = k_2 \begin{bmatrix} 1 \\ -2 \end{bmatrix}$$

We can also come up with an example where $\lambda_1 = \lambda_2$. For example:

$$\mathbf{M} = \begin{bmatrix} 0 & 1 \\ -1 & 2 \end{bmatrix}$$

$$\det(\mathbf{M} - \lambda \mathbf{I}) = 0$$

$$\det \left(\begin{bmatrix} -\lambda & 1 \\ -1 & 2 - \lambda \end{bmatrix} \right) = \lambda^2 - 2\lambda + 1 = (\lambda - 1)^2 = 0$$

Then we have $\lambda_1 = \lambda_2 = 1$.

Eigenvalues are **uniquely** determined (but the values can be repeated), while eigenvectors are **NOT**.

- ▶ Specifically, if ψ is an eigenvector, then $k\psi$ is as well, for any arbitrary constant real number k .
- ▶ If $\lambda_i = \lambda_{i+1}$, then $\psi_i + \psi_{i+1}$ will also be an eigenvector of eigenvalue λ_i . The eigenvectors of a given eigenvalue are only determined up to an orthogonal transformation.

$$\therefore (\lambda_i \mathbf{I} - \mathbf{M})\psi_i = (\lambda_i \mathbf{I} - \mathbf{M})\psi_{i+1} = 0$$

$$\therefore (\lambda_i \mathbf{I} - \mathbf{M})(\psi_i + \psi_{i+1}) = 0$$

Definition (1.3.2)

A matrix is positive definite if it is symmetric and all of its eigenvalues are positive. It is positive semidefinite if it is symmetric and all of its eigenvalues are nonnegative.

*When a **real** $n \times n$ matrix \mathbf{X} being positive definite: ^a*

$$\forall y \in \mathbb{R}^n, y^T \mathbf{X} y > 0$$

^a<https://mathworld.wolfram.com/PositiveDefiniteMatrix.html>

Fact (1.3.3)

The Laplacian matrix of a graph is positive semidefinite.

Proof (Fact 1.3.3)

Recall that previously we have that, for the Laplacian \mathbf{L}_G of (undirected) graph G , given a vector $\mathbf{x} \in \mathbb{R}^n$:

$$\mathbf{x}^T \mathbf{L}_G \mathbf{x} = \sum_{(a,b) \in E} w_{a,b} (\mathbf{x}(a) - \mathbf{x}(b))^2$$

when the weights $w_{a,b}$ are all non-negative, the value is non-negative as well.

In practice, we always number the eigenvalues of the Laplacian from **smallest** to **largest**.

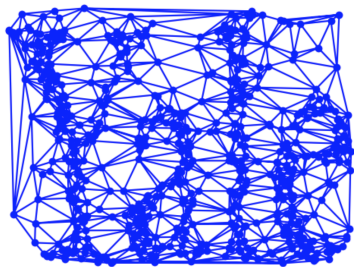
$$0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$$

We refer to $\lambda_2, \dots, \lambda_k$ (k is **small**) as low-frequency eigenvalues. λ_n is a high-frequency eigenvalue.

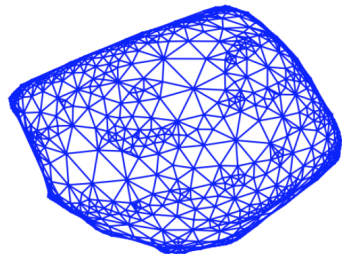
High and low frequency eigenmodes can be thought of as analogous to high and low frequency parts of the Fourier transform. ⁵

The second-smallest eigenvalue of the Laplacian matrix of a graph is zero ($\lambda_2 = 0$) iff the graph is disconnected. λ_2 is a measure of how well-connected the graph is. (See Chap 1.5.4 **The Fiedler Value**.)

⁵From a discussion on stackexchange.



(a) The original points sampled from Yale logo, with coordinates omitted and transformed into graph.



(b) Plot of vertices at $(\psi_2(a), \psi_3(a))$ coordinate.

Figure: An example showing the use of eigenvectors. More examples are listed in the textbook, Chap 1.

Eigenvalues and Optimization: The Courant-Fischer Theorem



One reason why we are interested in **eigenvalues** of matrices is that, they arise as the solution to natural **optimization** problems.

The formal statement of this is given by the **Courant-Fischer Theorem**. And this Theorem could be proved by the **Spectral Theorem**.

Theorem (2.0.1 Courant-Fischer Theorem)

Let \mathbf{M} be a symmetric matrix with eigenvalues $\mu_1 \geq \mu_2 \geq \cdots \geq \mu_n$. Then,

$$\mu_k = \max_{\substack{S \subseteq \mathbb{R}^n \\ \dim(S)=k}} \min_{\substack{\mathbf{x} \in S \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \min_{\substack{\mathcal{T} \subseteq \mathbb{R}^n \\ \dim(\mathcal{T})=n-k+1}} \max_{\substack{\mathbf{x} \in \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

where the maximization and minimization are over subspaces S and \mathcal{T} of \mathbb{R}^n .

Using the Spectral Theorem to prove the Courant-Fischer Theorem.

Theorem (1.3.1 The Spectral Theorem)

If \mathbf{M} is an n -by- n , real, symmetric matrix, then there exist real numbers $\lambda_1, \dots, \lambda_n$ and n mutually orthogonal unit vectors ψ_1, \dots, ψ_n and such that ψ_i is an eigenvector of \mathbf{M} of eigenvalue λ_i , for each i .

Main Steps:

- ▶ expanding a vector \mathbf{x} in the basis of eigenvectors of \mathbf{M}
- ▶ use the properties of eigenvalues and eigenvectors to prove it

$\mathbf{M} \in \mathbb{R}^{n \times n}$: a symmetric matrix, with eigenvalues $\mu_1 \geq \mu_2 \geq \cdots \geq \mu_n$. The corresponding **orthogonal** eigenvectors are $\psi_1, \psi_2, \dots, \psi_n$.

Then we may write $\mathbf{x} \in \mathbb{R}^n$ as:

$$\mathbf{x} = \sum_i c_i \psi_i, \quad c_i = \psi_i^T \mathbf{x}$$

Why \mathbf{x} can be expanded in this way?

Let Ψ be a matrix whose columns are $\{\psi_1, \psi_2, \dots, \psi_n\}$ — orthogonal vectors. By definition, Ψ is an orthogonal matrix.

$$\Psi\Psi^T = \Psi^T\Psi = I$$

Therefore we have:

$$\sum_i c_i \psi_i = \sum_i \psi_i c_i = \sum_i \psi_i \psi_i^T \mathbf{x} = \left(\sum_i \psi_i \psi_i^T \right) \mathbf{x} = \Psi\Psi^T \mathbf{x} = \mathbf{x}$$

and thus, since $\psi_i^T \psi_j = 1$ when $i = j$ and 0 otherwise,

$$\mathbf{x}^T \mathbf{x} = \left(\sum_i c_i \psi_i \right)^T \left(\sum_i c_i \psi_i \right) = \sum_{i,j} c_i^2 \psi_i^T \psi_j = \sum_{i=1}^n c_i^2$$

Let's revisit the theorem to prove (Now we have $\mathbf{x}^T \mathbf{x}$, to prove it we need to consider $\mathbf{x}^T \mathbf{M} \mathbf{x}$):

Theorem (2.0.1 Courant-Fischer Theorem)

Let \mathbf{M} be a symmetric matrix with eigenvalues $\mu_1 \geq \mu_2 \geq \cdots \geq \mu_n$. Then,

$$\mu_k = \max_{\substack{S \subseteq \mathbb{R}^n \\ \dim(S)=k}} \min_{\substack{\mathbf{x} \in S \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \max_{\substack{\mathcal{T} \subseteq \mathbb{R}^n \\ \dim(\mathcal{T})=n-k+1}} \min_{\substack{\mathbf{x} \in \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

where the maximization and minimization are over subspaces S and \mathcal{T} of \mathbb{R}^n .

In the textbook, Lemma 2.1.1 suggests that, in the previous example, for any $\mathbf{x} = \sum_i c_i \psi_i$:

$$\mathbf{x}^T \mathbf{M} \mathbf{x} = \sum_{i=1}^n c_i^2 \mu_i$$

Note that this time we **didn't** specify the value of c_i .

Again, $\psi_i^T \psi_j = 1$ when $i = j$ and 0 otherwise, also because $\mathbf{M}\psi_i = \mu_i\psi_i$,

$$\begin{aligned}\mathbf{x}^T \mathbf{M} \mathbf{x} &= \left(\sum_i c_i \psi_i \right)^T \mathbf{M} \left(\sum_i c_i \psi_i \right) \\ &= \left(\sum_i c_i \psi_i \right)^T \left(\sum_i c_i \mu_i \psi_i \right) \\ &= \sum_{i,j} c_i^2 \mu_i \psi_i^T \psi_j \\ &= \sum_i c_i^2 \mu_i\end{aligned}$$

Take a look again:

Theorem (2.0.1 Courant-Fischer Theorem)

Let \mathbf{M} be a symmetric matrix with eigenvalues $\mu_1 \geq \mu_2 \geq \cdots \geq \mu_n$. Then,

$$\mu_k = \max_{\substack{S \subseteq \mathbb{R}^n \\ \dim(S)=k}} \min_{\substack{\mathbf{x} \in S \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \max_{\substack{T \subseteq \mathbb{R}^n \\ \dim(T)=n-k+1}} \min_{\substack{\mathbf{x} \in T \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

where the maximization and minimization are over subspaces S and T of \mathbb{R}^n .

We need the value of $\frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$. In particular, we care about μ_k and subspace \mathcal{S} where $\dim(\mathcal{S}) = k$. Also recall that we put $\{\mu_i\}_{i=1}^n$ in the **non-increasing** order.

$$\mathbf{x} = \sum_i^k c_i \boldsymbol{\psi}_i, \quad c_i = \boldsymbol{\psi}_i^T \mathbf{x}$$

$$\frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \frac{\sum_i^k c_i^2 \mu_i}{\sum_i^k c_i^2} \geq \frac{\sum_i^k c_i^2 \mu_k}{\sum_i^k c_i^2} = \mu_k$$

Therefore,

$$\min_{\substack{\mathbf{x} \in \mathcal{S} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \geq \mu_k$$

To prove the theorem, we also need to show that for all subspace $\mathcal{S} \subseteq \mathbb{R}^n$ where $\dim(\mathcal{S}) = k$,

$$\min_{\substack{\mathbf{x} \in \mathcal{S} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \leq \mu_k$$

For this part we bring up the subspace \mathcal{T} of dimension $n - k + 1$, whose basis vectors are ψ_k, \dots, ψ_n . Similarly, for $\mathbf{x} \in \mathcal{T}$, we have:

$$\max_{\substack{\mathbf{x} \in \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \frac{\sum_k^n c_i^2 \mu_i}{\sum_k^n c_i^2} \leq \frac{\sum_k^n c_i^2 \mu_k}{\sum_k^n c_i^2} = \mu_k$$

Every subspace \mathcal{S} of dimension k has an intersection with \mathcal{T} (dimension $n - k + 1$), the intersection has dimension at least 1 ($((n - k + 1) + k = n + 1)$).

$$\min_{\substack{\mathbf{x} \in \mathcal{S} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \leq \min_{\substack{\mathbf{x} \in \mathcal{S} \cap \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \leq \max_{\substack{\mathbf{x} \in \mathcal{S} \cap \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \leq \max_{\substack{\mathbf{x} \in \mathcal{T} \\ \mathbf{x} \neq \mathbf{0}}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} = \mu_k$$

The theorem is proved this way.

We prove the Spectral Theorem in a form that is almost identical to Courant-Fischer.

Main Steps:

- ▶ showing that the **Rayleigh quotient** and eigenvectors, eigenvalues have certain relation, starting from μ_1 ;
- ▶ use the conclusion in the first step to prove that a vector is an eigenvector, prove the Spectral Theorem by generalizing this characterization to **all** of the eigenvalues of \mathbf{M}

The **Rayleigh quotient** of a vector \mathbf{x} with respect to a matrix \mathbf{M} is defined to be:

$$\frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

The Rayleigh quotient of an eigenvector is its corresponding eigenvalue: if $\mathbf{M}\boldsymbol{\psi} = \mu\boldsymbol{\psi}$, then (by default, $\boldsymbol{\psi} \neq \mathbf{0}$)

$$\frac{\boldsymbol{\psi}^T \mathbf{M} \boldsymbol{\psi}}{\boldsymbol{\psi}^T \boldsymbol{\psi}} = \frac{\boldsymbol{\psi}^T (\mathbf{M} \boldsymbol{\psi})}{\boldsymbol{\psi}^T \boldsymbol{\psi}} = \frac{\boldsymbol{\psi}^T (\mu \boldsymbol{\psi})}{\boldsymbol{\psi}^T \boldsymbol{\psi}} = \frac{\mu \boldsymbol{\psi}^T \boldsymbol{\psi}}{\boldsymbol{\psi}^T \boldsymbol{\psi}} = \mu$$

The first step is to prove the following theorem:

Theorem (2.2.1 (Rayleigh quotient and eigenvectors))

Let \mathbf{M} be a symmetric matrix and let vector $\mathbf{x} \neq \mathbf{0}$ maximize the **Rayleigh quotient** with respect to \mathbf{M} :

$$\frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

Then, $\mathbf{M}\mathbf{x} = \mu_1 \mathbf{x}$, where μ_1 is the largest eigenvalue of \mathbf{M} . Conversely, the minimum is achieved by eigenvectors of the smallest eigenvalue of \mathbf{M} .

Observe that:

- ▶ the Rayleigh quotient is homogeneous, maximum is achieved;
- ▶ it suffices to consider unit vectors \mathbf{x}
- ▶ the set of unit vectors is a closed and compact set where the maximum is achieved on

Recall that: a function at its maximum and minimum has gradient $\mathbf{0}$ (zero vector).

We can compute the gradient of the Rayleigh quotient.

$$\nabla \mathbf{x}^T \mathbf{x} = 2\mathbf{x} \quad \mathbf{x}^T \mathbf{M} \mathbf{x} = 2\mathbf{M} \mathbf{x}$$

also recall the derivative rule:

$$\left(\frac{f}{g}\right)' = \frac{gf' - fg'}{g^2}$$

$$\nabla \left(\frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \right) = \frac{(\mathbf{x}^T \mathbf{x})(2\mathbf{M} \mathbf{x}) - (\mathbf{x}^T \mathbf{M} \mathbf{x})(2\mathbf{x})}{(\mathbf{x}^T \mathbf{x})^2}, \quad \mathbf{x} \neq \mathbf{0}$$

when it is $\mathbf{0}$, $(\mathbf{x}^T \mathbf{x})\mathbf{M} \mathbf{x} = (\mathbf{x}^T \mathbf{M} \mathbf{x})\mathbf{x}$, $\mathbf{M} = \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}} \mathbf{x}$.

Recall that: *the Rayleigh quotient of an eigenvector is its corresponding eigenvalue.* The above equation holds iff \mathbf{x} is an eigenvector of \mathbf{M} . Proved.

Theorem (2.2.2 (almost identical to the CF Theorem))

Let M be an n -dimensional real symmetric matrix. There exist numbers μ_1, \dots, μ_n and orthonormal vectors ψ_1, \dots, ψ_n such that $M\psi_i = \mu_i\psi_i$. Moreover,

$$\psi_1 \in \arg \max_{\|\mathbf{x}\|=1} \mathbf{x}^T M \mathbf{x}$$

and for $2 \leq i \leq n$,

$$\psi_i \in \arg \max_{\substack{\|\mathbf{x}\|=1 \\ \mathbf{x}^T \psi_j = 0, j < i}} \mathbf{x}^T M \mathbf{x},$$

$$\text{similarly, } \psi_i \in \arg \min_{\substack{\|\mathbf{x}\|=1 \\ \mathbf{x}^T \psi_j = 0, j > i}} \mathbf{x}^T M \mathbf{x}$$

To start with, we want to reduce to the case of positive definite matrices. In order to do that, we first modify \mathbf{M} a bit.

$$\mu_n = \min_{\mathbf{x}} \frac{\mathbf{x}^T \mathbf{M} \mathbf{x}}{\mathbf{x}^T \mathbf{x}}$$

we know μ_n exists from Theorem 2.2.1 we've just proved. Now we consider:

$$\widetilde{\mathbf{M}} = \mathbf{M} + (1 - \mu_n) \mathbf{I}$$

For $\forall \mathbf{x}$ such that $\|\mathbf{x}\| = 1$, we have:

$$\mathbf{x}^T \widetilde{\mathbf{M}} \mathbf{x} = \mathbf{x}^T \mathbf{M} \mathbf{x} + 1 - \mu_n = 1 + (\mathbf{x}^T \mathbf{M} \mathbf{x} - \min_{\mathbf{x}} \mathbf{x}^T \mathbf{M} \mathbf{x}) \geq 1$$

Therefore $\widetilde{\mathbf{M}}$ is positive definite.

Besides,

$$\widetilde{\mathbf{M}}\mathbf{x} = \mathbf{M}\mathbf{x} + (1 - \mu_n)\mathbf{x}$$

For $\forall \psi, \mu$ where $\mathbf{M}\psi = \mu\psi$,

$$\widetilde{\mathbf{M}}\psi = \mathbf{M}\psi + (1 - \mu_n)\psi = (\mu + 1 - \mu_n)\psi$$

thus $\widetilde{\mathbf{M}}$ and \mathbf{M} have the same eigenvectors.

Thus it suffices to prove the theorem for positive definite matrices. In other words, we treat \mathbf{M} as if it is positive definite.

We proceed by induction on k . We construct ψ_{i+1} base on eigenvalues ψ_1, \dots, ψ_k satisfying:

$$\psi_i \in \arg \max_{\substack{\|\mathbf{x}\|=1 \\ \mathbf{x}^T \psi_j = 0, j < i}} \mathbf{x}^T \mathbf{M} \mathbf{x}$$

And define:

$$\mathbf{M}_k = \mathbf{M} - \sum_{i=1}^k \mu_i \psi_i \psi_i^T$$

For $j \leq k$ we have (because all the previous eigenvectors are all orthogonal to each other):

$$\mathbf{M}_k \psi_j = \mathbf{M} \psi_j - \sum_{i=1}^k \mu_i \psi_i \psi_i^T \psi_j = \mu_j \psi_j - \mu_j \psi_j = \mathbf{0}$$

Hence, for vector \mathbf{x} that are orthogonal to ψ_1, \dots, ψ_k ,

$$\mathbf{M}\mathbf{x} = \mathbf{M}_k\mathbf{x} + \sum_{i=1}^k \mu_i \psi_i \psi_i^T \mathbf{x} = \mathbf{M}_k\mathbf{x}, \quad \mathbf{x}^T \mathbf{M}\mathbf{x} = \mathbf{x}^T \mathbf{M}_k\mathbf{x}$$

and,

$$\arg \max_{\substack{\|\mathbf{x}\|=1 \\ \mathbf{x}^T \psi_j = 0, j < i}} \mathbf{x}^T \mathbf{M}\mathbf{x} \leq \arg \max_{\|\mathbf{x}\|=1} \mathbf{x}^T \mathbf{M}_k\mathbf{x}$$

For convenience we define $\mathbf{y} = \arg \max_{\|\mathbf{x}\|=1} \mathbf{x}^T \mathbf{M}_k\mathbf{x}$. From Theorem 2.2.1 we know that \mathbf{y} is an eigenvector of \mathbf{M}_k . Let's say that the corresponding eigenvalue is μ . \mathbf{M}_k and \mathbf{M} have the same eigenvectors, thus \mathbf{y} is an eigenvector of \mathbf{M} .

Now we will prove that we can set $\psi_{k+1} = \mathbf{y}$ and $\mu_{k+1} = \mu$.

We prove it by showing \mathbf{y} must be orthogonal to each ψ_1, \dots, ψ_k .

$$\tilde{\mathbf{y}} = \mathbf{y} - \sum_{i=1}^k \psi_i(\psi_i^T \mathbf{y})$$

is the projection of \mathbf{y} orthogonal to ψ_1, \dots, ψ_k . Since $\mathbf{M}_k \psi_j = \mathbf{0}$ for $j \leq k$,

$$\tilde{\mathbf{y}}^T \mathbf{M}_k \tilde{\mathbf{y}} = \mathbf{y}^T \mathbf{M}_k \mathbf{y} = \mathbf{y}^T \mathbf{M} \mathbf{y}$$

If \mathbf{y} is not orthogonal to ψ_1, \dots, ψ_k , some $\psi_i^T \mathbf{y} \neq \mathbf{0}$, then $\|\tilde{\mathbf{y}}\| < \|\mathbf{y}\|$. Because we assume positive definite of \mathbf{M} , there comes a contradiction.

$$\tilde{\mathbf{y}}^T \mathbf{M}_k \tilde{\mathbf{y}} = \tilde{\mathbf{y}}^T \mathbf{M} \tilde{\mathbf{y}} > 0$$

and also note that $\|\tilde{\mathbf{y}}\| < \|\mathbf{y}\|$ (previous conclusion), for normalized $\hat{\mathbf{y}}, \hat{\mathbf{y}} = \tilde{\mathbf{y}}/\|\tilde{\mathbf{y}}\|$, and \mathbf{y} was an **unit** vector,

$$\begin{aligned}\hat{\mathbf{y}}^T \mathbf{M} \hat{\mathbf{y}} &= \hat{\mathbf{y}}^T \mathbf{M}_k \hat{\mathbf{y}} = \frac{\tilde{\mathbf{y}}^T \mathbf{M}_k \tilde{\mathbf{y}}}{\|\tilde{\mathbf{y}}\|^2} \\ &> \frac{\tilde{\mathbf{y}}^T \mathbf{M}_k \tilde{\mathbf{y}}}{\|\mathbf{y}\|^2} = \tilde{\mathbf{y}}^T \mathbf{M}_k \tilde{\mathbf{y}} = \mathbf{y}^T \mathbf{M}_k \mathbf{y} = \mathbf{y}^T \mathbf{M} \mathbf{y}\end{aligned}$$

There's a conflict with \mathbf{y} 's definition:

$$\mathbf{y} = \arg \max_{\|\mathbf{x}\|=1} \mathbf{x}^T \mathbf{M}_k \mathbf{x}$$

$\therefore \mathbf{y}$ must be orthogonal to ψ_1, \dots, ψ_k . Proved.

The Laplacian and Graph Drawing



Chapter 3 shows that Laplacian should reveal a lot about the structure of graphs, although not always guaranteed to work.

It mentions Hall's (Kenneth M. Hall) work a lot of times:
An r -dimensional quadratic placement algorithm

The idea of drawing graphs using eigenvectors demonstrated in Section 1.5.1 was suggested by Hall in 1970.

Recall that weighted undirected graph $G = (V, E, w)$, with positive weight $w : E \rightarrow \mathbb{R}^+$, is defined this way:

$$\mathbf{L}_G \stackrel{\text{def}}{=} \mathbf{D}_G - \mathbf{M}_G, \quad \mathbf{D}_G = \sum_b w_{a,b}$$

where \mathbf{D}_G is the diffusion matrix, \mathbf{M}_G is the adjacency matrix.

Given a vector $\mathbf{x} \in \mathbb{R}^n$,

$$\mathbf{x}^T \mathbf{L}_G \mathbf{x} = \sum_{(a,b) \in E} w_{a,b} (\mathbf{x}(a) - \mathbf{x}(b))^2$$

Here are some of the very interesting properties of a graph that we would like to prove.

- ▶ If and only if the graph is connected, there are only one eigenvalue equals to zero.
- ▶ When mapping each vertex to a real number, so as to create their coordinates for drawing, orthogonal eigenvectors are useful.

Lemma

Let $G = (V, E)$ be a graph, and let $0 = \lambda_1 \leq \lambda_2 \leq \dots \leq \lambda_n$ be the eigenvalues of its Laplacian matrix, \mathbf{L} . Then, $\lambda_2 > 0$ if and only if G is connected.

First of all, there exists eigenvalue $\mathbf{0}$, because the all-one vector $\mathbb{1}$ satisfies:

$$\mathbf{L}\mathbb{1} = \mathbf{0}$$

To prove, if we view the Laplacian $\mathbf{L} = \mathbf{D} - \mathbf{M}$ as an operator ($\mathbf{D} = \sum_{(a,b) \in E} w_{a,b}$), for each \mathbf{x} we have its a^{th} entry of $\mathbf{L}\mathbf{x}$ being:

$$(\mathbf{L}\mathbf{x})(a) = d(a)\mathbf{x}(a) - \sum_{(a,b) \in E} w_{a,b}\mathbf{x}(b) = \sum_{(a,b) \in E} w_{a,b}(\mathbf{x}(a) - \mathbf{x}(b))$$

It infers that $\mathbb{1}$ is an eigenvector corresponds to eigenvalue 0. Therefore, $\lambda_1 = 0$.

Next, we show that $\lambda_2 = 0$ if G is disconnected.

If G is disconnected, then we can split it into two graphs G_1 and G_2 . Because we can safely reorder the vertices of a graph, we can have:

$$\mathbf{L} = \begin{bmatrix} \mathbf{L}_{G_1} & \mathbf{0} \\ \mathbf{0} & \mathbf{L}_{G_2} \end{bmatrix}$$

It has at least 2 orthogonal eigenvectors of eigenvalue zero:

$$\begin{bmatrix} \mathbf{0} \\ \mathbf{1} \end{bmatrix}, \text{ and } \begin{bmatrix} \mathbf{1} \\ \mathbf{0} \end{bmatrix}$$

On the other hand, for a eigenvector ψ of eigenvalue 0, $\mathbf{L}\psi = \mathbf{0}$,

$$\psi^T \mathbf{L} \psi = \sum_{(a,b) \in E} w_{a,b} (\psi(a) - \psi(b))^2 = 0$$

For every pair of vertices (a, b) connected by an edge, we have $\psi(a) = \psi(b)$. In a connected graph, all vertices are directly or indirectly connected, and thus ψ must be a constant vector.

Contradiction found.

Therefore, G must be disconnected when $\lambda_2 = 0$.

Hall's idea on graph drawing suggests that we choose the first coordinates of the n vertices as $\mathbf{x} \in \mathbb{R}^n$ that minimizes:

$$\mathbf{x}^T \mathbf{L} \mathbf{x} = \sum_{(a,b) \in E} w_{a,b} (\mathbf{x}(a) - \mathbf{x}(b))^2 = 0$$

To avoid degenerating to $\mathbf{0}$, we have restriction:

$$\|\mathbf{x}\|^2 = \sum_{a \in V} \mathbf{x}(a)^2 = 1$$

To avoid degenerating to $\mathbb{1}/\sqrt{n}$, Hall suggested another constraint:

$$\mathbb{1}^T \mathbf{x} = \sum_{a \in V} \mathbf{x}(a) = 0$$

When there are multiple sets of coordinates, say \mathbf{x} and \mathbf{y} ; we require $\mathbf{x}^T \mathbf{y} = 0$, to avoid cases such as $\mathbf{x} = \mathbf{y} = \psi_2$.

Theorem (3.2.1)

Let \mathbf{L} be a Laplacian matrix and let $\mathbf{x}_1, \dots, \mathbf{x}_k$ be orthonormal vectors that are all orthogonal to $\mathbf{1}$. Then

$$\sum_{i=1}^k \mathbf{x}_i^T \mathbf{L} \mathbf{x}_i \geq \sum_{i=2}^{k+1} \lambda_i$$

and this inequality is tight only when $\mathbf{x}^T \boldsymbol{\psi}_j = 0$ for all j such that $\lambda_j \geq \lambda_{k+1}$. λ_i are the eigenvalues, the graph G is an undirected connected graph.

We can order λ such that:

$$0 = \lambda_1 \leq \lambda_2 \leq \cdots \leq \lambda_n$$

As is proved before, $\lambda_1 = 0$ and because G is connected, ψ_1 is a constant vector.

Let $\mathbf{x}_{k+1} \dots \mathbf{x}_n$ be vectors such that $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_n$ is an orthogonal **basis**. It is done by choosing $\mathbf{x}_{k+1} \dots \mathbf{x}_n$ to be an orthogonal basis of the space orthogonal to $\mathbf{x}_1, \mathbf{x}_2, \dots, \mathbf{x}_k$. Because they are orthogonal basis, (think of orthogonal matrix)

$$\sum_{j=1}^n (\psi_j^T \mathbf{x}_i)^2 = \sum_{j=1}^n (\mathbf{x}_i^T \psi_j)^2 = 1, \quad i = 1, 2, \dots, n$$

Because of that $\boldsymbol{\psi}_1^T \mathbf{x}_i \propto \mathbb{1}^T \mathbf{x}_i = 0$, and that $\sum_{j=1}^n (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = 1$,

$$\sum_{j=2}^n (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = 1$$

Previously, $\mathbf{x}^T \mathbf{M} \mathbf{x} = \sum_i c_i^2 \mu_i$, $c_i = \boldsymbol{\psi}_i^T \mathbf{x}$, $\mathbf{x} = \sum_i c_i \boldsymbol{\psi}_i$. Here,

$$\begin{aligned} \mathbf{x}_i^T \mathbf{L} \mathbf{x}_i &= \sum_{j=2}^n \lambda_j (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = \lambda_{k+1} + \sum_{j=2}^n (\lambda_j - \lambda_{k+1}) (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \\ &\geq \lambda_{k+1} + \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1}) (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \end{aligned}$$

It is tight only when $\boldsymbol{\psi}_j^T \mathbf{x}_i = 0$ for $\lambda_j \geq \lambda_{k+1}$.

$$\lambda_{k+1} + \sum_{j=2}^n (\lambda_j - \lambda_{k+1})(\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \geq \lambda_{k+1} + \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1})(\boldsymbol{\psi}_j^T \mathbf{x}_i)^2$$

Quick proof of when the above inequality is tight:

$$\begin{aligned} \lambda_{k+1} + \sum_{j=2}^n (\lambda_j - \lambda_{k+1})(\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 &= \lambda_{k+1} + \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1})(\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \\ &\quad + \sum_{j=k+2}^n (\lambda_j - \lambda_{k+1})(\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = 0 \end{aligned}$$

That is $\boldsymbol{\psi}_j^T \mathbf{x}_i = 0$ for $j > k + 1$. When $j > k + 1$, $\lambda_j \geq \lambda_{k+1}$.

To prove the Theorem 3.2.1, we sum up over i :

$$\begin{aligned}\sum_i^k \mathbf{x}_i^T \mathbf{L} \mathbf{x}_i &\geq k\lambda_{k+1} + \sum_{i=1}^k \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1}) (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \\ &= k\lambda_{k+1} + \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1}) \sum_{i=1}^k (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \\ &\geq k\lambda_{k+1} + \sum_{j=2}^{k+1} (\lambda_j - \lambda_{k+1}) = \sum_{j=2}^{k+1} \lambda_j\end{aligned}$$

because: $\lambda_j - \lambda_{k+1} \leq 0$, and, $\sum_{i=1}^k (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \leq \sum_{i=1}^n (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = 1$.