

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



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## Introduction

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# Introduction

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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*<sup>1</sup>.

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  $\mathbf{W}_G \mathbf{p}$ .

<sup>1</sup>Regular 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  $\delta_a$  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: <sup>2</sup>

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

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).

<sup>2</sup>Note 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  $\psi$  is an eigenvector of a matrix  $\mathbf{M}$  with eigenvalue  $\lambda$  if:

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

$\lambda$  is an eigenvalue if and only if  $\lambda\mathbf{I} - \mathbf{M}$  is a singular matrix ( $\because \det(\lambda\mathbf{I} - \mathbf{M}) = 0$ ). 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 “ $\lambda$  is an eigenvalue if and only if  $\lambda\mathbf{I} - \mathbf{M}$  is a singular matrix”: <sup>3</sup>

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

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

is a homogeneous linear system for  $\psi$ , 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.

<sup>3</sup>[https://www-users.math.umn.edu/~olver/num\\_/lnv.pdf](https://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  $\psi_1, \dots, \psi_n$  and such that  $\psi_i$  is an eigenvector of  $\mathbf{M}$  of eigenvalue  $\lambda_i$ , 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. <sup>4</sup>

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

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

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

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  $\psi_1$  and  $\psi_2$ , 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  $\psi$  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  $\lambda_i$ . 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: <sup>a</sup>*

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

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<sup>a</sup><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**.

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

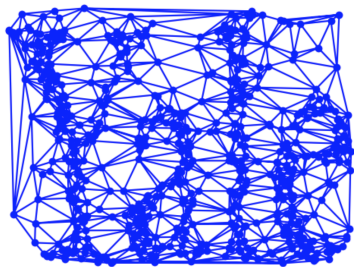
We refer to  $\lambda_2, \dots, \lambda_k$  ( $k$  is **small**) as low-frequency eigenvalues.  $\lambda_n$  is a high-frequency eigenvalue.

*High and low frequency eigenmodes can be thought of as analogous to high and low frequency parts of the Fourier transform.*<sup>5</sup>

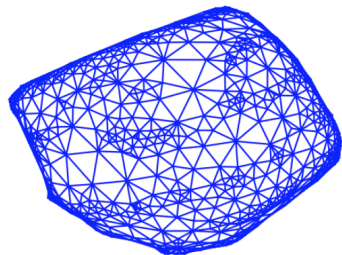
The second-smallest eigenvalue of the Laplacian matrix of a graph is zero ( $\lambda_2 = 0$ ) iff the graph is disconnected.  $\lambda_2$  is a measure of how well-connected the graph is. (See Chap 1.5.4 **The Fiedler Value**.)

<sup>5</sup>From 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

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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,

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

where the maximization and minimization are over subspaces  $\mathcal{S}$  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  $\psi_1, \dots, \psi_n$  and such that  $\psi_i$  is an eigenvector of  $\mathbf{M}$  of eigenvalue  $\lambda_i$ , 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  $\Psi$  be a matrix whose columns are  $\{\psi_1, \psi_2, \dots, \psi_n\}$  — orthogonal vectors. By definition,  $\Psi$  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  $\mu_k$  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  $\psi_k, \dots, \psi_n$ . 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  $\mu_1$ ;
- ▶ 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  $\mu_1$  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  $\mu_1, \dots, \mu_n$  and orthonormal vectors  $\psi_1, \dots, \psi_n$  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  $\mu_n$  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  $\forall \boldsymbol{\psi}, \mu$  where  $\mathbf{M}\boldsymbol{\psi} = \mu\boldsymbol{\psi}$ ,

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

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  $\psi_{i+1}$  base on eigenvalues  $\psi_1, \dots, \psi_k$  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  $\psi_1, \dots, \psi_k$ ,

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

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  $\mu$ .  $\mathbf{M}_k$  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  $\psi_1, \dots, \psi_k$ .

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

is the projection of  $\mathbf{y}$  orthogonal to  $\psi_1, \dots, \psi_k$ . 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  $\psi_1, \dots, \psi_k$ , 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  $\psi_1, \dots, \psi_k$ . Proved.

# The Laplacian and Graph Drawing

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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  $\psi$  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  $\psi$  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  $1/\sqrt{n}$ , Hall suggested another constraint:

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

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  $\lambda_j \geq \lambda_{k+1}$ .  $\lambda_i$  are the eigenvalues, the graph  $G$  is an undirected connected graph.

We can order  $\lambda$  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,  $\psi_1$  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,  $\sum_{i=1}^k (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 \leq \sum_{i=1}^n (\boldsymbol{\psi}_j^T \mathbf{x}_i)^2 = 1$ .