

1 Spectral Graph Theory

We assume the graph has n vertices and m edges. The vertex set is V and the edge set is E .

1.1 Courant-Fischer Theorem

1. **Eigenvalue version:** Let A be a symmetric matrix with eigenvalues $\lambda_1 \leq \dots \leq \lambda_n$, then

$$\lambda_i = \min_{\substack{\text{subspace } W \subseteq \mathbb{R}^n \\ \dim(W)=i}} \max_{\substack{x \in W \\ x \neq 0}} \frac{x^\top A x}{x^\top x} = \max_{\substack{\text{subspace } W \subseteq \mathbb{R}^n \\ \dim(W)=n+1-i}} \min_{\substack{x \in W \\ x \neq 0}} \frac{x^\top A x}{x^\top x}.$$

2. **Eigenbasis version:** Let A be a symmetric matrix with eigenvalues $\lambda_1 \leq \dots \leq \lambda_n$ and corresponding orthonormal eigenvectors x_1, \dots, x_n , then

$$\lambda_i = \min_{\substack{x \perp x_1, \dots, x_{i-1} \\ x \neq 0}} \frac{x^\top A x}{x^\top x} = \max_{\substack{x \perp x_{i+1}, \dots, x_n \\ x \neq 0}} \frac{x^\top A x}{x^\top x}.$$

Note that we also have $\lambda_i = \frac{x_i^\top A x_i}{x_i^\top x_i}$.

Applying Courant-Fischer theorem, we have $\lambda_2 x^\top x \leq x^\top L x \leq \lambda_n x^\top x$ for all $x \perp \mathbf{1}$, as $\mathbf{1}$ is the eigenvector of L corresponding to eigenvalue 0. For connected graphs, $\lambda_2 > 0$.

1.2 PSD Order (Loewner Order)

Defined only for symmetric matrices: $A \leq B$ iff for all $x \in \mathbb{R}^n$, we have $x^\top A x \leq x^\top B x$. We also define $G \leq H$ for two graphs G and H iff $L_G \leq L_H$. We always have $G \geq H$ if H is a subgraph of G .

Properties:

1. If $A \leq B$ and $B \leq C$, then $A \leq C$.
2. If $A \leq B$, then $A + C \leq B + C$ for any symmetric C .
3. If $A \leq B$ and $C \leq D$, then $A + C \leq B + D$.
4. If $A > 0$ and $\alpha \geq 1$, then $\frac{1}{\alpha} A \leq A \leq \alpha A$.
5. If $A \leq B$, then $\lambda_i(A) \leq \lambda_i(B)$ for all i . Proof by Courant-Fischer theorem. The converse is not true.
6. For any matrix C , if $A \leq B$, then $C^\top A C \leq C^\top B C$.
7. If $0 \leq A \leq B$, then $B^{-1} \leq A^{-1}$.

1.3 Bounding the λ_2 and λ_n

1.3.1 Test Vector Method

Since $\lambda_2 \leq \frac{y^\top L y}{y^\top y}$ for any $y \perp \mathbf{1}$, we can upper bound the λ_2 by any **test vector** y . Similarly, we can lower bound the λ_n by test vectors by $\lambda_n \geq \frac{y^\top L y}{y^\top y}$.

1. For a complete graph K_n , $L = nI - \mathbf{1}\mathbf{1}^\top$ and for any $x \perp \mathbf{1}$ we have $x^\top L x = n x^\top x$. Therefore, $\lambda_2(K_n) = \dots = \lambda_n(K_n) = n$ and any $x \perp \mathbf{1}$ is an eigenvector.
2. For a path graph P_n , let $x(i) = n + 1 - 2i$ be the test vector which satisfies $x \perp \mathbf{1}$, we get $\lambda_2(P_n) \leq \frac{12}{n^2}$. Let $x(1) = -1$, $x(n) = 1$ and $x(i) = 0$ for other i to be the test vector, we get $\lambda_n(P_n) \geq 1$.
3. For a complete binary tree T_n (depth equals zero for a single root), let $x(i) = 0$ for all non-leaf nodes, $x(i) = -1$ for

even-numbered leaf nodes and $x(i) = 1$ for odd-numbered leaf nodes be the test vector, we get $\lambda_n(T_n) \geq 1$. Let $x(1) = 0$, $x(i) = 1$ for the left subtree of the root and $x(i) = -1$ for the right subtree of the root be the test vector, we get $\lambda_2(T_n) \leq \frac{2}{n-1}$.

1.3.2 Consequences of PSD Order

Since $x^\top (D - A)x = \sum_{(u,v)} w(u,v)(x(u) - x(v))^2 \geq 0$ and $x^\top (D + A)x = \sum_{(u,v)} w(u,v)(x(u) + x(v))^2 \geq 0$, we have $D \geq A$ and $D \geq -A$. In addition, we have $D \leq (\max D_{i,i})I$. Therefore, we have $L = D - A \leq 2D \leq (2 \max D_{i,i})I$, which implies $\lambda_n \leq 2 \max D_{i,i}$ for any graph. For unit-weight graphs, this means $\lambda_n \leq 2 \max \text{degree}(v)$. The bound is tight for a single-edge graph.

To get lower bounds of $\lambda_2(H)$, we first establish $f(n)H \geq G$ for some G with known lower bounds on $\lambda_2(G)$. Usually $G = K_n$ because $\lambda_2(K_n) = n$. Then it follows that $\lambda_2(H) \geq \lambda_2(G)/f(n)$.

1. **Path Graph P_n :** Let $G_{i,j}$ denote a unit-weight graph consisting of one edge (i, j) and P_n be the path graph connecting 1 and n . Then $(n-1)P_n \geq G_{1,n}$. Proof follows from applying Cauchy-Schwartz for $\delta_i := x(i+1) - x(i)$. For weighted paths, we have $G_{1,n} \leq \left(\sum_{i=1}^{n-1} \frac{1}{w_i}\right) \sum_{i=1}^{n-1} w_i G_{i,i+1}$. Applying path inequality, we have $K_n = \sum_{i < j} G_{i,j} \leq \sum_{i < j} (j-i)P_{i,j} \leq \sum_{i < j} (j-i)P_n \leq n^3 P_n$, which implies $\lambda_2(P_n) \geq \lambda_2(K_n)/n^3 = 1/n^2$.
2. **Any unit-weight graph G :** Define the diameter D of a graph G to be the maximum length of the shortest paths between any two nodes. Let $G_{i,j}^s$ be the shortest path from i to j . Applying path inequality, we have $K_n = \sum_{i < j} G_{i,j} \leq \sum_{i < j} D G_{i,j}^s \leq \sum_{i < j} D G \leq n^2 D G$, which implies $\lambda_2(G) \geq \frac{1}{nD}$.
3. **Complete Binary Tree T_n :** Define G_e be the single-edge graph with edge e , and $T_{i,j}$ be the unique path between i and j . Applying the weighted path inequality, we have $K_n = \sum_{i < j} G_{i,j} \leq \sum_{i < j} \left(\left(\sum_{e \in T_{i,j}} \frac{1}{w(e)} \right) \left(\sum_{e \in T_{i,j}} w(e) G_e \right) \right) \leq \left(\max_{i < j} \sum_{e \in T_{i,j}} \frac{1}{w(e)} \right) \left(\sum_{i < j} \sum_{e \in T_{i,j}} w(e) G_e \right)$. For e connecting level i and $i+1$ for $i \in [d-1]$, we set $w(e) = 2^i$. Then $\max_{i < j} \sum_{e \in T_{i,j}} \frac{1}{w(e)} \leq 4$. Since the number of occurrence of e in $T_{i,j}$ for any $i < j$ is upper bounded by $n^2 2^{-i}$, we have $\sum_{i < j} \sum_{e \in T_{i,j}} w(e) G_e \leq \sum_e w(e) n^2 2^{-i} G_e = \sum_e n^2 G_e = n^2 T_n$. Therefore, $K_n \leq 4n^2 T_n$, which implies $\lambda_2(T_n) \geq \frac{1}{4n}$.

2 Conductance

Definitions:

1. **Conductance of a vertex subset:** Given $\emptyset \subset S \subset V$, the conductance $\phi(S) := \phi(S) = \frac{|E(S, V \setminus S)|}{\min\{\text{vol}(S), \text{vol}(V \setminus S)\}}$, where $\text{vol}(S) := \sum_{v \in S} \text{degree}(v)$. Define $\mathbf{1}_S$ to be the n -dimensional vector with only 1 for the vertices of S and 0 for the vertices of $V \setminus S$. Assuming $\text{vol}(S) \leq \text{vol}(V)/2$, thus $|E(V, V \setminus S)| =$

$\sum_{(u,v) \in E} (\mathbf{1}_S(u) - \mathbf{1}_S(v))^2 = \mathbf{1}_S^\top L \mathbf{1}_S$ and $\text{vol}(S) = \mathbf{1}_S^\top D \mathbf{1}_S$. Then $\phi(S) = \frac{\mathbf{1}_S^\top L \mathbf{1}_S}{\mathbf{1}_S^\top D \mathbf{1}_S}$.

2. **Conductance of a graph:** The conductance $\phi(G) := \min_{\emptyset \subset S \subset V} \phi(S) = \min_{\substack{\emptyset \subset S \subset V \\ \text{vol}(S) \leq \text{vol}(V)/2}} \phi(S)$.
3. **ϕ -expander:** For any $\phi \in (0, 1]$, we call a graph G to be a ϕ -expander if $\phi(G) \geq \phi$.
4. **ϕ -expander decomposition of quality q :** A partition $\{X_i\}$ of the vertex set V is called a ϕ -expander decomposition of quality q if (1) each induced graph $G[X_i]$ is a ϕ -expander, and (2.i) #edges not contained in any $G[X_i]$ is at most $q \cdot \phi \cdot m$. The second condition is equivalent to (2.ii) The partition removes at most $q \cdot \phi \cdot m$ edges.
5. **Normalized Laplacian:** We define the *normalized Laplacian* to be $N := D^{-1/2} L D^{-1/2}$. N is still PSD, with first eigenvalue equals 0 associated with eigenvector $D^{1/2} \mathbf{1}$. By Courant-Fischer theorem, $\lambda_2(N) = \min_{x \perp D^{1/2} \mathbf{1}} \frac{x^\top N x}{x^\top x} = \min_{z \perp d} \frac{z^\top L z}{z^\top D z}$.

2.1 Cheeger's Inequality

Notice that the $\lambda_2(N)$ has similar forms to $\phi(G)$. Cheeger's inequality aims to bound $\phi(G)$ by $\lambda_2(N)$.

Cheeger's Inequality: $\frac{\lambda_2(N)}{2} \leq \phi(G) \leq \sqrt{2 \lambda_2(N)}$.

The lower bound is proved by restricting the minimum in $\lambda_2(N)$ to be $z_S = \mathbf{1}_S - \alpha \mathbf{1}$ for some α such that $z_S \perp d$. The upper bound is proved by constructing S for any $z \perp d$ such that $\frac{\mathbf{1}_S^\top L \mathbf{1}_S}{\mathbf{1}_S^\top D \mathbf{1}_S} \leq \sqrt{2 \frac{z^\top L z}{z^\top D z}}$.

3 Random Walks on a Graph

A **random walk on a graph** G is a Markov Chain with transition probability $\mathbb{P}(v_{t+1} = v \mid v_t = u) = w(u, v)/d(u)$ iff $(u, v) \in E$ and 0 otherwise. The transition matrix is thus $W = A D^{-1} = I - D^{-1/2} N D^{-1/2}$ and $p_t = W^t p_0$. Define $\pi = \frac{d}{1^\top d}$, thus $\pi = W \pi$ for any G , so every G has a stationary distribution.

3.1 Lazy Random Walks

A **lazy random walk on a graph** G is a random walk, but has half probability to not move for every step. Assuming that G is connected, the lazy random walk guarantees ergodicity of the Markov Chain, and thus convergence to the stationary distribution. The transition matrix is $\tilde{W} = \frac{1}{2}(I + W) = I - \frac{1}{2} D^{-1/2} N D^{-1/2}$.

Relation between lazy random walk and normalized Laplacian: For the i -th eigenvalue v_i of N associated with eigenvector ψ_i , the \tilde{W} has an eigenvalue $1 - \frac{1}{2} v_i$ associated with eigenvector $D^{1/2} \psi_i$. Since $0 \leq L \leq 2D$, we have $0 \leq N \leq 2I$ and thus $0 \leq \lambda_i(N) \leq 2$. Therefore, we conclude that all eigenvalues of $\tilde{W} \in [0, 1]$.

Dynamics of lazy random walk: Expanding the starting distribution p_0 by the eigenvectors of \tilde{W} , we have for some $\{\alpha_i\}$ that $p_0 = \sum_{i=1}^n \alpha_i D^{1/2} \psi_i$. Therefore, we have $p_t = \tilde{W}^t p_0 = \sum_{i=1}^n \alpha_i (1 - \frac{1}{2} v_i)^t D^{1/2} \psi_i \rightarrow \alpha_1 D^{1/2} \psi_1$ as $v_1 = 0$ and

$v_i > 0$ for $i \neq 1$. Since $\psi_1 \propto D^{1/2} \mathbf{1}$, we have $\psi_1 = \frac{d^{1/2}}{(\mathbf{1}^\top d)^{1/2}}$, thus $\alpha_1 = \psi_1^\top D^{-1/2} p_0 = \frac{\mathbf{1}^\top p_0}{(\mathbf{1}^\top d)^{1/2}} = \frac{1}{(\mathbf{1}^\top d)^{1/2}}$ and $\alpha_1 D^{1/2} \psi_1 = \pi$, which implies $p_t \rightarrow \pi$, the stationary distribution.

Rate of Convergence: For any unit-weight connected graph G and any starting distribution p_0 , we have $\|p_t - \pi\|_\infty \leq e^{-v_2 t/2} \sqrt{n}$. Therefore, a larger v_2 and smaller vertex set means faster convergence, and the convergence rate is exponential.

3.2 Hitting Time

The expected hitting time from a to s is defined by $\mathbb{E}H_{a,s}$, where $H_{a,s} = \text{argmin}_t \{v_t = s \mid v_0 = a\}$. We want $\mathbb{E}H_{a,s}$ for all vertices a and denote the vector as h , e.g., $h(s) = 0$.

By one-step analysis, we have $h(a) = 1 + \sum_{(a,b) \in E} \frac{w(a,b)}{d(a)} h(b) = 1 + \mathbf{1}_a^\top W^\top h$, and thus $\mathbf{1} = \mathbf{1}_a^\top (\mathbf{I} - W^\top) h$. Combining the equation for all vertices except s , we have $\mathbf{1} - \alpha \mathbf{1}_s = (\mathbf{I} - W^\top) h$, where α represents the extra freedom from the $n-1$ equations. Multiplying both side by D , we get $d - \alpha d(s) \mathbf{1}_s = (D - A)h = Lh$, which only have solution when $d - \alpha d(s) \mathbf{1}_s \perp \mathbf{1}$. Therefore, $\alpha = \|d\|_1 / d(s)$.

To summarize, by solving $Lh = d - \|d\|_1 \mathbf{1}$, we can get the expected hitting time from all vertices to s . Note that the solution has one extra freedom because $\dim(\ker(L)) = 1$, and the correct expected hitting time is $h - h(s) \mathbf{1}$ to enforce the constraint that $h(s) = 1$. The equation can be solved in $\tilde{O}(m)$.

4 Pseudo-Inverse and Effective Resistance

Given a Laplacian L , its (Moore-Penrose) pseudo inverse is defined to be either of the two equivalents:

1. A matrix L^+ that is (1) symmetric, (2) $L^+ v = 0$ for $v \in \ker(L)$, and (3) $L^+ L v = L L^+ v = v$ for $v \in \ker(L)$.
2. Let λ_i, v_i be the i -th eigenvalue and eigenvector. Then $L^+ = \sum_{\lambda_i \neq 0} \lambda_i^{-1} v_i v_i^\top$.

Property:

- Assume $M = XYX^\top$, where X is real and invertible, and Y is real and symmetric. Let Π_M be the orthogonal projection to the image of M . Then $M^+ = \Pi_M (X^\top)^{-1} Y^+ X^{-1} \Pi_M$.
- For symmetric L , $\Pi_L := \sum_{\lambda_i \neq 0} v_i v_i^\top = L^{+1/2} L L^{+1/2} = L^+ L = L L^+$ is the orthogonal projection to the image of L , i.e., $\Pi_L v = 0$ for any $v \in \ker(L)$ and $\Pi_L v = v$ for any $v \in \text{im}(L)$. For connected G , $\Pi_L = \mathbf{I} - \frac{1}{n} \mathbf{1} \mathbf{1}^\top$.

The effective resistance between vertex a and b is defined to be the cost (energy lost) to routing one unit (of positive electric charge) from a to b : $R_{\text{eff}}(a, b) = \min_{Bf = \mathbf{1}_b - \mathbf{1}_a} f^\top R f = \tilde{f}^\top R \tilde{f}$, where \tilde{f} is the electric flow. Let \tilde{x} be the electric voltages, we also have $L\tilde{x} = \mathbf{1}_b - \mathbf{1}_a$, and thus $R_{\text{eff}}(a, b) = \tilde{x}^\top L \tilde{x} = (\mathbf{1}_b - \mathbf{1}_a)^\top L^+ (\mathbf{1}_b - \mathbf{1}_a) = \|L^{+1/2} (\mathbf{1}_b - \mathbf{1}_a)\|_2^2$.

Effective Resistance is a distance defined on the vertex pairs, i.e. $R_{\text{eff}}(a, c) \leq R_{\text{eff}}(a, b) + R_{\text{eff}}(b, c)$.

5 Gaussian Elimination for Laplacian

5.1 Optimization View

Solving $Lx = d$ is equivalent to solving $\text{argmin}_x -d^\top x + \frac{1}{2} x^\top L x$. By iteratively optimize over x_i , we get a series of similar optimizations. The final optimization is straightforward, then we can back substitute to get x .

5.2 Additive View

Given an invertible square lower/upper triangular matrix M , we can solve $Mx = d$ by back substitution in $O(\text{nnz}(M))$, where $\text{nnz}(M)$ means the number of non-zeros in M . Therefore, if we know the **Cholesky decomposition** $L = MM^\top$ (requires $O(n^3)$), then we can solve $Lx = d = M(M^\top x)$ by (1) solving $My = d$ then (2) solving $M^\top x = y$ in $O(\text{nnz}(M))$.

However, the Laplacian is non-invertible, leading to one diagonal of M equals 0. Therefore, we need to play a trick. Define \hat{M} equals M but has value 1 for the zero diagonal, and \hat{D} be a diagonal matrix that has value 0 at the zero diagonal of M and 1 otherwise. Then $L = \hat{M} \hat{D} \hat{M}^\top$, and each \hat{M} is now invertible. Since $\hat{D}^+ = \hat{D}$, we can find a special solution of $Lx = d$ by (1) solving $\hat{M}z = d$, (2) computing $y = \hat{D}z$, and (3) solving $\hat{M}^\top x = y$. The solution space is obtained by adding a subspace spanned by $\mathbf{1}$.

6 Approximating a Dense Graph in the Spectral Domain

6.1 Concentration of Random Matrices

1. **Chernoff Bound for Bounded independent variables:** Suppose $\{X_i \in \mathbb{R}\}$ are independent random variables and $0 \leq X_i \leq R$. Let $X = \sum_i X_i$ and $\mu = \mathbb{E}X$. Then for any $0 < \epsilon \leq 1$, we have $\mathbb{P}(X \geq (1 + \epsilon)\mu) \leq \exp(-\frac{\epsilon^2 \mu}{4R})$ and $\mathbb{P}(X \leq (1 - \epsilon)\mu) \leq \exp(-\frac{\epsilon^2 \mu}{4R})$.
2. **Bernstein Bound for independent, zero-mean and bounded variables:** Suppose $\{X_i \in \mathbb{R}\}$ are independent, zero-mean random variables and $|X_i| \leq R$. Let $X = \sum_i X_i$, and $\sigma^2 = \text{Var}(X)$. Then for any $t > 0$, we have $\mathbb{P}(|X| \geq t) \leq 2 \exp(-\frac{t^2}{2Rt + 4\sigma^2})$. The proof is similar to Chernoff bound. (1) $\mathbb{P}(X \geq t) = \mathbb{P}(\exp(\theta X) \geq \exp(\theta t)) \leq \exp(-\theta t) \mathbb{E}(\exp(\theta X))$, (2) upper bound $\mathbb{E}(\exp(\theta X)) \leq \exp(\theta^2 \sigma^2)$ given $\theta \in (0, \frac{1}{R}]$, which allows $\exp(\theta X_i) \leq 1 + \theta X_i + (\theta X_i)^2$, and (3) take the minimum among $\theta \in (0, \frac{1}{R}]$.
3. **Bernstein Bound for independent, zero-mean and bounded symmetric matrices:** Suppose $\{X_i \in \mathbb{R}^{n \times n}\}$ are independent, zero-mean, symmetric random matrices and $\|X_i\| \leq R$, where $\|\cdot\|$ is the spectral norm (the largest singular value). Let $X = \sum_i X_i$, and $\sigma^2 = \|\sum_{i=1}^n \mathbb{E}X_i^2\|$. Then $\mathbb{P}(\|X\| \geq t) \leq 2n \exp(-\frac{t^2}{2Rt + 4\sigma^2})$.

6.2 Matrix Functions

Given a real-valued function $f : \mathbb{R} \rightarrow \mathbb{R}$ and a symmetric matrix A with eigen-decomposition $A = V \Lambda V^\top$, we define

$f(A) = V f(\Lambda) V^\top$. This is compatible to the Taylor expansion $f(x) = \sum_i \alpha_i x^i$, as $f(A) = \sum_i \alpha_i A^i = V (\sum_i \alpha_i f(\Lambda)) V^\top = V f(\Lambda) V^\top$.

Monotonicity: Given $f : \mathcal{D} \rightarrow \mathcal{C}$ and partial orders $\leq_{\mathcal{C}}$ and $\leq_{\mathcal{D}}$, we call f is monotonically increasing w.r.t. these orders iff for all $d_1 \leq_{\mathcal{D}} d_2 \in \mathcal{D}$ we have $f(d_1) \leq_{\mathcal{C}} f(d_2)$. For matrix functions, we use the PSD order as the ordering.

Property:

- If the scalar function f is monotonically increasing, then the matrix function $X \rightarrow \text{Tr}(f(X))$ is monotonically increasing.
- $\log(\cdot)$ is monotonically increasing.
- The matrix function $(\cdot)^2$ and $\exp(\cdot)$ is **not** monotone.
- $\exp(A) \leq \mathbf{I} + A + A^2$ for $\|A\| \leq 1$.
- $\log(\mathbf{I} + A) \leq A$ for $A > -\mathbf{I}$.
- (Lieb's theorem): $f(A) := \text{Tr}(\exp(H + \log(A)))$ for some symmetric H is concave in the domain of PSD matrices. This is in particular useful with Markov theorem.

6.3 Approximating a Dense Graph by Sparse Graphs

Given PD matrices A, B and $\epsilon > 0$, we say $A \approx_{\epsilon} B$ iff $\frac{1}{1+\epsilon} A \leq B \leq (1+\epsilon)A$. If $L_G \approx_{\epsilon} L_{\tilde{G}}$ and $|\tilde{E}| \ll |E|$, we call \tilde{G} a spectral sparsifier of G .

Properties:

- Define $c_G(T) := \sum_{e \in E \cap (T \times V \setminus T)} w(e)$ to be the value of the cut $(T, V \setminus T)$. If $L_G \approx_{\epsilon} L_{\tilde{G}}$, then for all $T \subset V$, we have $\frac{1}{1+\epsilon} c_G(T) \leq c_{\tilde{G}}(T) \leq (1+\epsilon) c_G(T)$. The proof is by noticing $c_G(T) = \mathbf{1}_T^\top L_G \mathbf{1}_T$.
 - $L \approx_{\epsilon} \tilde{L} \Leftrightarrow \Pi_L \approx_{\epsilon} L^{+1/2} \tilde{L} L^{+1/2}$, as $A \leq B$ implies $C^\top A C \leq C^\top B C$ for any $C \in \mathbb{R}^{n \times n}$.
 - For $\epsilon \leq 1$, if $\|\Pi_L - L^{+1/2} \tilde{L} L^{+1/2}\| \leq \epsilon/2$, then $\Pi_L \approx_{\epsilon} L^{+1/2} \tilde{L} L^{+1/2}$.
- Theorem:** Consider a connected graph $G = (V, E, w)$, with $n = |V|$. For any $0 < \epsilon < 1$ and $0 < \delta < 1$, there exist sampling probabilities p_e for each edge $e \in E$ s.t. if we include each edge e in \tilde{E} independently with probability p_e and set its weight $\tilde{w}(e) = \frac{1}{p_e} w(e)$, then with probability at least $1 - \delta$ the graph $\tilde{G} = (V, \tilde{E}, \tilde{w})$ satisfies $L_G \approx_{\epsilon} L_{\tilde{G}}$ and $|\tilde{E}| \leq O(n \epsilon^{-2} \log(n/\delta))$. The proof uses Bernstein bounds to prove the concentration of the constructed random graph.

7 Solving Laplacian Linear Equations Approximately

Idea: solving Laplacian linear equations requires $O(m)$, which is expensive when the graph is dense. By approximating the Laplacian, we can get an approximation of the solution quickly.

Given PSD matrix M and $d \in \text{im}(M)$, let $Mx^* = d$. We say that \tilde{x} is an ϵ -approximate solution to $Mx = d$ iff $\|\tilde{x} - x^*\|_M^2 \leq \epsilon \|x^*\|_M^2$, where $\|x\|_M^2 = x^\top M x$. Note that any solution to $Mx = d$ has the same $\|\cdot\|_M^2$, as they differ by a vector in the kernel of M .

Theorem: Given a Laplacian L of a weighted undirected graph $G = (V, E, w)$ with $|E| = m$ and $|V| = n$ and a demand vector $d \in \mathbb{R}^V$, we can find \tilde{x} that is an ϵ -approximate solution to

$Lx = d$, using an algorithm that takes time $O(m \log^c n \log(1/\epsilon))$ for some fixed constant c and succeeds with probability $1 - 1/n^{10}$. Note that without known Cholesky decomposition in advance, the exact solution requires $O(n^3)$ and $m \leq n^2/2$.