

Approximating the Exponential, the Lanczos Method and an $\tilde{O}(m)$ -Time Spectral Algorithm for Balanced Separator

Lorenzo Orecchia, MIT

Sushant Sachdeva, Princeton University

Nisheeth Vishnoi, MSR India

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BALANCED GRAPH PARTITIONING AND SPECTRAL METHODS

For an undirected unweighted instance graph $G=(V,E)$ with $|V|=n$ and $|E|=m$, the conductance of a cut $S \subseteq V$ is defined as

$$\phi(S) = \frac{|E(S, \bar{S})|}{\min\{\text{vol}(S), \text{vol}(\bar{S})\}}$$

where $\text{vol}(S)$ is the total degree in set S . A cut S is b -balanced if $\text{vol}(S) \geq b \cdot \text{vol}(V)$.

b -BALANCED CUT PROBLEM: Given graph G , parameter $\gamma \in (0,1)$ and balance $b \in (0,1/2)$, does G contain a b -balanced cut of conductance at most γ ?

This problem is NP-hard. However, approximation algorithms exist:

Algorithm	Method	Distinguishes $\geq \gamma$ and	Running Time
Iterative Eigenvector	Spectral	$O(\sqrt{\gamma})$	$\tilde{O}(n^2)$
[Leighton, Rao '88]	Flow	$O(\gamma \log n)$	$\tilde{O}(m^{\frac{3}{2}})$ [AK '07, OSVV '08]
[Arora, Rao, Vazirani '04]	SDP (Flow + Spectral)	$O(\gamma \sqrt{\log n})$	$\tilde{O}(m^{\frac{3}{2}})$ [Sherman '09]
[Madry '10]	SDP (Flow + Spectral)	$\gamma \text{polylog}(n)$	$\tilde{O}(m^{1+\epsilon})$

We use "spectral methods" to refer to algorithms that explore the graph by performing matrix-vector multiplications involving the graph Laplacian L . Such algorithms detect low-conductance cuts by exploiting the connection between the mixing of random walks in the graph and the cut structure of G .

Spectral methods for finding balanced cuts perform well in many applications and have fast running times and optimized implementations that make them a popular choice among practitioners. The following are the theoretical guarantees of some spectral algorithms for this problem.

Algorithm	Method	Distinguishes $\geq \gamma$ and	Time
Iterative Eigenvector	Eigenvector	$O(\sqrt{\gamma})$	$\tilde{O}(n^2)$
[Andersen, Peres '09]	Local Random Walks	$O(\sqrt{\gamma \log n})$	$\tilde{O}(m/\sqrt{\gamma})$
[Orecchia, Vishnoi '10]	SDP	$O(\sqrt{\gamma})$	$\tilde{O}(m/\gamma)$

OUR THEOREM: We give an algorithm that either outputs an $\Omega(b)$ -balanced cut $S \subseteq V$ such that $\phi(S) \leq O(\sqrt{\gamma})$, or outputs a certificate that no b -balanced cut of conductance γ exists. The algorithm runs in time $O(m \text{poly}(\log n))$.

TECHNICAL COMPONENTS:

- 1) SDP primal-dual iterative algorithm with a simple random walk interpretation
- 2) Novel analysis of Lanczos methods for computing heat-kernel vectors

2

DETECTING AND REMOVING UNBALANCED CUTS

Unbalanced cuts of low conductance are obstacles to detecting balanced cuts.

Spectral methods are targeted towards finding low-conductance cuts, regardless of how balanced they are. For this reason, spectral methods may have to find and remove all unbalanced cuts of low conductance before finding a balanced cut.

The algorithmic challenge is to detect and remove unbalanced cuts of low conductance quickly.

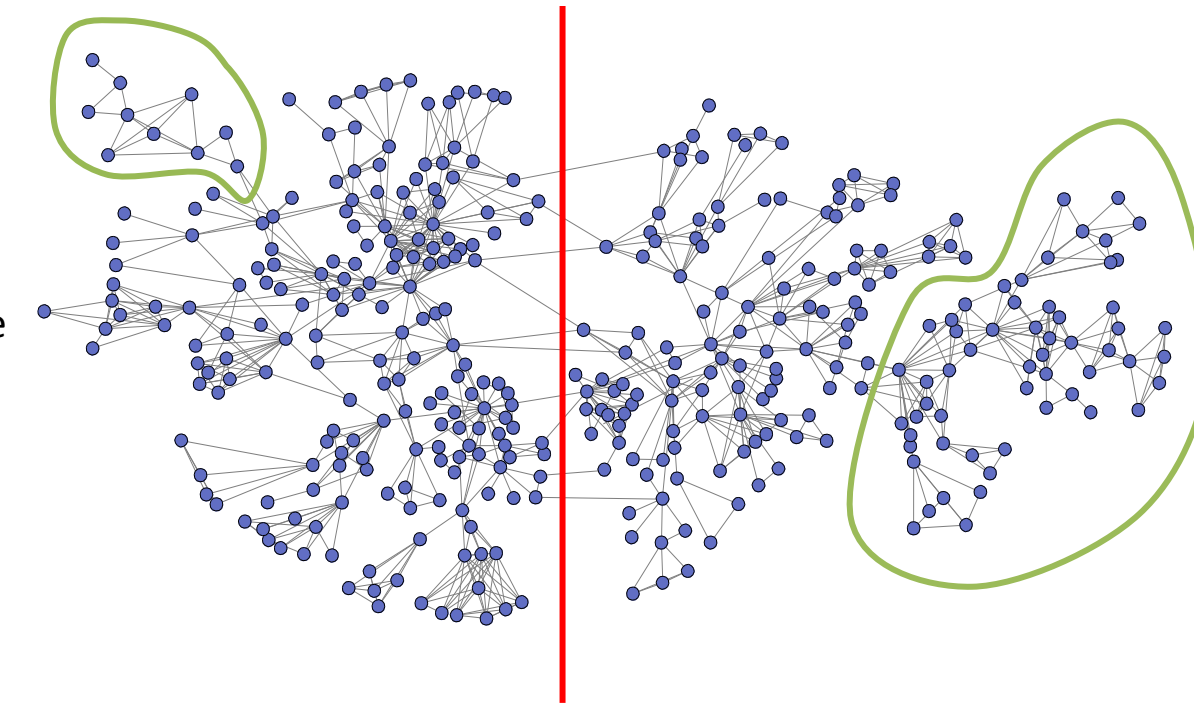


Figure: Balanced cuts and unbalanced cuts, both of low conductance, in a citation graph.

Iterative Eigenvector Approach

Set $G_t = G$. For $t=1, \dots, n$ do:

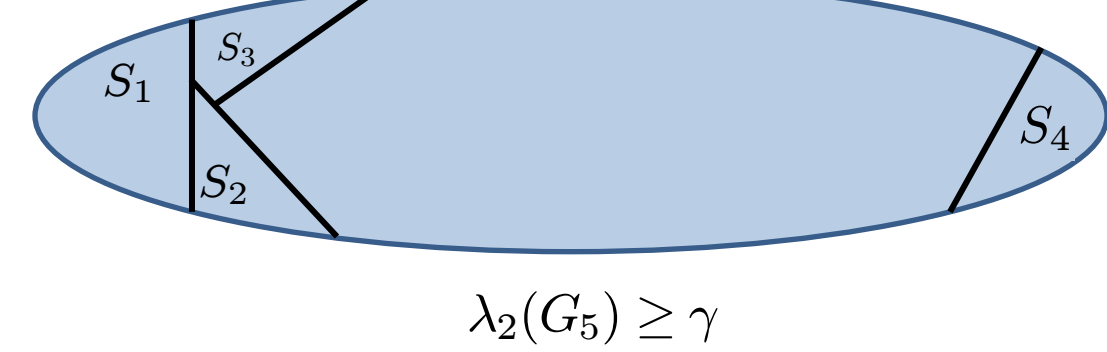
- Compute the slowest mixing eigenvector of G_t and corresponding Laplacian eigenvalue λ_t .

- If $\lambda_t \geq \gamma$, output NO. Otherwise, sweep eigenvector to find cut S_t such that $\phi(S_t) \leq O(\sqrt{\gamma})$.

- If U_{S_t} is $(b/2)$ -balanced, output S_t .

- Otherwise, let G_{t+1} be the graph induced by G_t on $V - S_t$ with self-loops replacing the edges going to S_t . Recurse on G_{t+1} .

Example:

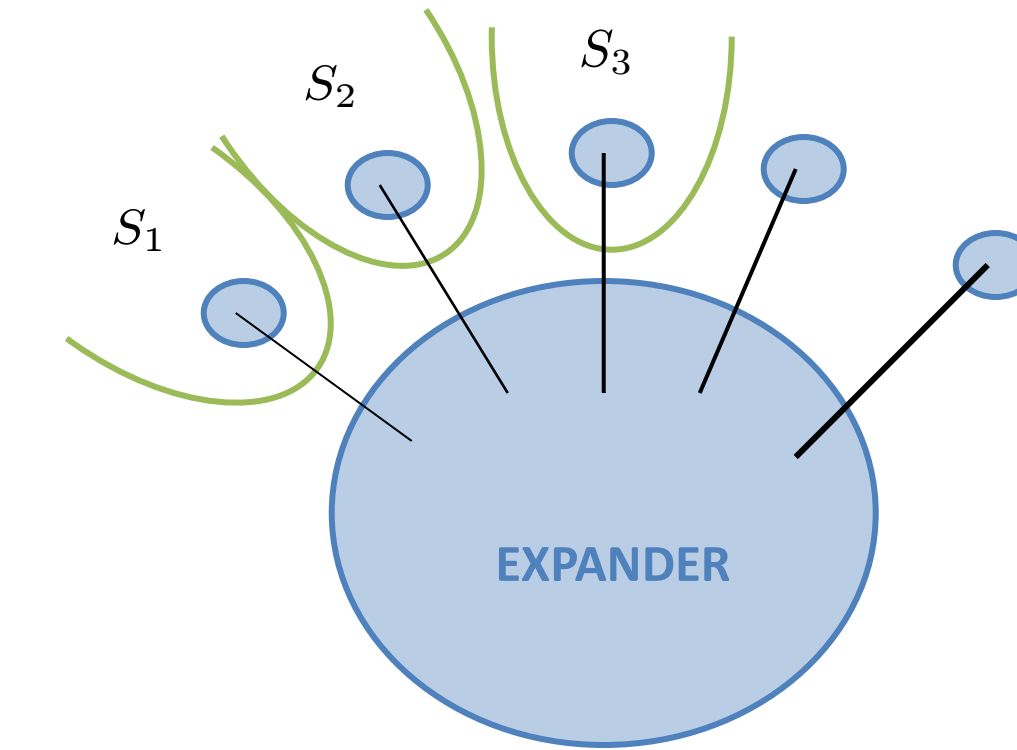


Quadratic Running Time

Worst-case instance: $\sim (n)$ iterations.

Each iteration takes time $\sim (n)$, yielding a total running time of $\sim (n^2)$.

NB: It is not possible to argue that $\lambda_2(G_t)$ grows significantly at every iteration.



3

OUR RANDOM-WALK APPROACH FOR FINDING UNBALANCED CUTS

The eigenvector is unstable, as it may capture only a small cut and be oblivious to rest of the graph. Hence, we switch to a more stable distribution over low eigenvectors, represented by the transition matrix of a random walk that has converged to have most of its norm over eigenvectors with eigenvalue at most $O(\gamma)$. For simplicity, take G to be regular. The graph Laplacian is $L = I - W$.

IDEA BEHIND OUR ALGORITHM:

Replace eigenvector by multidimensional embedding of heat-kernel random-walk.

$$P_t = e^{-\tau L(G_t)} \quad \tau = \log n / \gamma$$

Denote the embedding $\{v_i\}$ given by $v_i = P e_i$ with $P = P_t$.

MIXING:

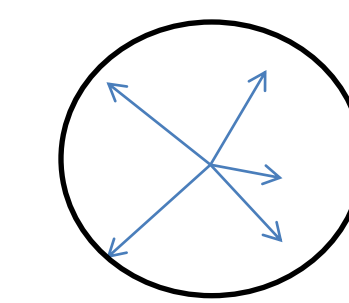
Define the deviation from stationary for a set $S \subseteq V$ for process P :

$$\Psi(P, S) = \sum_{i \in S} \|v_i - \bar{1}/n\|^2$$

$\Psi(P, V)$ measures the total deviation from stationary, which is also the variance of the geometric embedding $\{v_i\}$.

ROUNDING. Three possible cases:

WALK HAS MIXED



All vectors are short. The total deviation is:

$$\Psi(P_t, V) \leq \frac{1}{\text{poly}(n)}$$

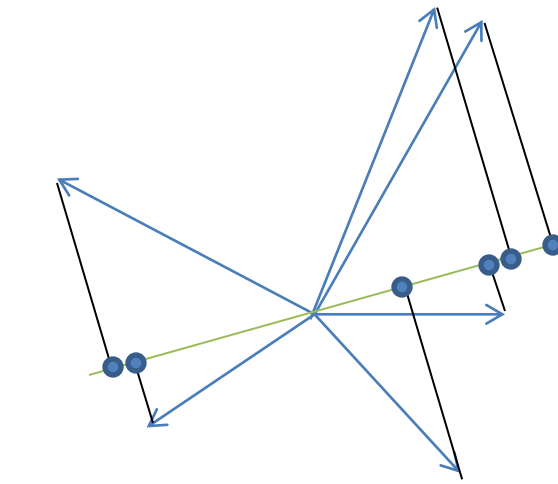
This yields a certificate that

$$\lambda_2(G_t) \geq \gamma,$$

Which we turn into a NO certificate.

WALK HAS NOT MIXED AND BALANCED CUT IS FOUND

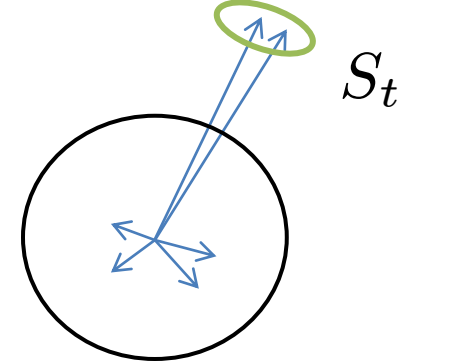
We take $O(\log n)$ random projections and find the best $\Omega(b)$ -balanced sweep cut S of the resulting vectors.



If $\phi(S) \leq O(\sqrt{\gamma})$, we output S and answer YES.

WALK HAS NOT MIXED AND NO BALANCED CUT IS FOUND

Some vectors must be very long, i.e. walk converges very poorly from some vertices.



We can find S_t with $\phi(S_t) \leq O(\sqrt{\gamma})$.

Moreover,

$$\Psi(P_t, S_t) \geq \frac{1}{2} \Psi(P_t, V).$$

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HOW WE UPDATE THE GRAPH AND OUR MIXING ANALYSIS

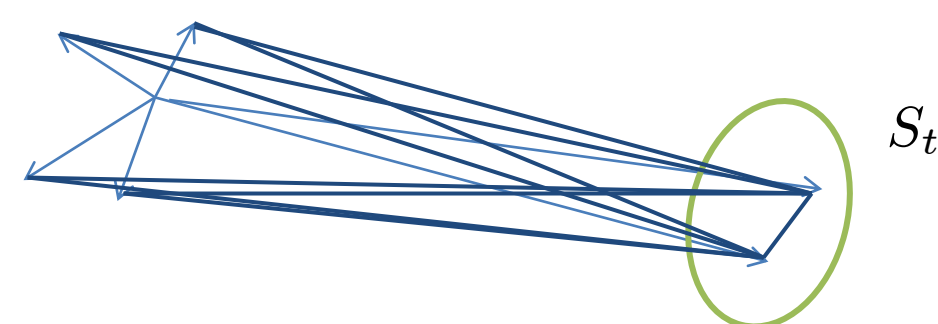
If a low-conductance balanced cut is found, we perform a **soft removal** of this cut by **modifying the current random walk** as follows:

$$P_{t+1} = e^{-\tau \left(L + \gamma \sum_{j=1}^{t-1} \sum_{i \in S_j} L(\text{Star}_i) \right)}$$

where Star_i is the **star graph** rooted at vertex i .

The transition rate from S_t to all vertices increases, making the process converge faster to its stationary.

Increased Transition Rates



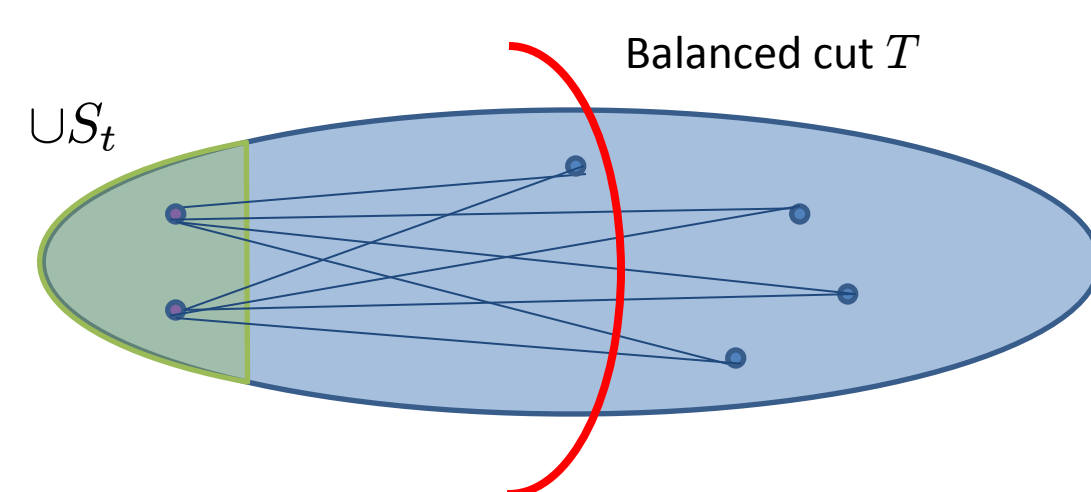
MIXING ANALYSIS: Using properties of the heat-kernel random walk, we show that the mixing improves significantly:

$$\Psi(P_{t+1}, V) \leq \Psi(P_t, V) - \frac{1}{2} \Psi(P_t, S_t) \leq \frac{3}{4} \Psi(P_t, V)$$

After $T=O(\log n)$ iterations, if no low-conductance $\Omega(b)$ -balanced cut is found, the following holds:

$$\Psi(P_T, V) \leq \frac{1}{\text{poly}(n)}$$

We show that we can turn this fact into a **NO** certificate for the b -Balanced Cut problem:



$$\phi(T) \geq \gamma - \gamma \frac{|U_{S_t}|}{|T|} \geq \gamma - \gamma \frac{b/2}{b} \geq \frac{\gamma}{2}$$

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HOW TO COMPUTE OUR RANDOM-WALK VECTORS

GOAL: For symmetric diagonally-dominant A and vector u , with sparsity m , compute vector v such that

$$\|e^{-A} u - v\| \leq \delta \quad \text{in time } \tilde{O}(m \cdot \text{polylog}(\|A\|) \cdot \text{polylog}(1/\delta))$$

In our case, $A = \frac{\log n}{\gamma} L$, so that $\|A\| = \tilde{O}(\frac{1}{\gamma})$ and the running time is $\tilde{O}(m)$.

ITERATIVE APPROACHES:

- **Taylor Series Approximation:** requires $\Theta(\|A\| \log(1/\delta))$ terms and yields a running time of $\tilde{O}(m/\gamma)$.

- **Direct Lanczos Method:** requires $\Theta(\sqrt{\|A\|} \text{polylog}(1/\delta))$ iterations. The running time is $\tilde{O}(m/\sqrt{\gamma})$.

- **Our Algorithm** relies on a linear-system solver for A . We obtain the following running times:

- Using Spielman-Teng solver: $\tilde{O}((m+n) \cdot \log(2+\|A\|) \cdot \text{polylog}(1/\delta))$, yielding $\tilde{O}(m)$

- Using Conjugate Gradient: $\tilde{O}((m+n) \cdot \sqrt{1+\|A\|} \cdot \log(2+\|A\|) \cdot \text{polylog}(1/\delta))$, yielding $\tilde{O}(m/\sqrt{\gamma})$

OUR APPROACH exploits the speed of the linear-system solver by using it to perform the **inverse iteration**. We speed up convergence by applying **Lanczos method**.

Review of Lanczos Method

IDEA: Perform k matrix-vector multiplications to obtain the subspace

$$R_k = \text{Span}\{u, Au, A^2u, A^3u, A^4u, \dots, A^k u\}$$

Compute an orthonormal basis Q_k for R_k and let T_k be A restricted to R_k :

$$T_k = Q_k A Q_k^T$$

As k grows, T_k becomes a better approximation to A . If a function f is close to a polynomial p of degree k , i.e.

$$\|p(x) - f(x)\| \leq \delta \quad \forall x \in \text{Spectrum}(A)$$

then

$$\|Q_k f(T_k) Q_k^T u - f(A)u\| \leq 2\delta \|u\|$$

6

APPLYING LANCZOS METHOD TO THE INVERSE ITERATION

A direct application of Lanczos method does not meet our goal, but applying Lanczos to the inverse iteration yield our result.

DIRECT LANCZOS: There exists a polynomial p such that

$$\sup_{x \in (0, \|A\|)} |p(x) - e^{-x}| \leq \frac{1}{\text{poly}(n)}$$

and p has degree $\tilde{O}(\sqrt{\|A\|})$.

IMPLICATION: $\tilde{O}(\sqrt{\|A\|}) = \tilde{O}(1/\sqrt{\gamma})$ iterations are sufficient.

LOWER BOUND: This bound is tight: $\tilde{O}(\sqrt{\|A\|})$ iterations may be necessary.

IDEA: Apply Lanczos Method to $B = (I + \frac{A}{k})^{-1}$, i.e. compute subspace

$$R_k = \text{Span}\{u, Bu, B^2u, B^3u, B^4u, \dots, B^k u\}$$

NB: We can use the linear-system solvers to compute, for any vector x ,

$$Bx = (I + \frac{A}{k})^{-1} x$$

APPROXIMATION: In fact, the quality of the k -approximation depends on the existence of a good approximation to the exponential function by rational functions, a larger class of functions than polynomials.

THEOREM (Saff, Schonhage, Varga '75): There exists a polynomial p such that

$$\sup_{x \in (0, \|A\|)} \left| \frac{p(x)}{(1+x/k)^k} - e^{-x} \right| \leq \delta$$

and p has degree $\tilde{O}(\log(2+\|A\|) \cdot \text{polylog}(1/\delta))$.

STRONG ERROR GUARANTEE