

COMPSCI 514: ALGORITHMS FOR DATA SCIENCE

Andrew McGregor

Lecture 21

Computing the SVD/eigendecomposition

- Efficient algorithms for SVD/eigendecomposition.
- High level: a glimpse into fast methods for linear algebraic computation, which are workhorses behind data science.

We have talked about the eigendecomposition and SVD as ways to compress data, to embed entities like words and documents, to compress/cluster non-linearly separable data.

How efficient are these techniques? Can they be run on massive datasets?

Power Method: The most fundamental iterative method for approximate SVD/eigendecomposition. Applies to computing $k = 1$ eigenvectors, but can be generalized to larger k .

Goal: Given symmetric $\mathbf{A} \in \mathbb{R}^{d \times d}$, with eigendecomposition $\mathbf{A} = \mathbf{V}\mathbf{\Lambda}\mathbf{V}^T$, find \vec{z} which is an approximation to the top eigenvector \vec{v}_1 of \mathbf{A} .

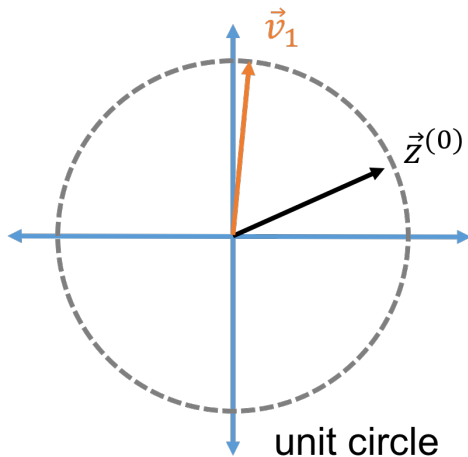
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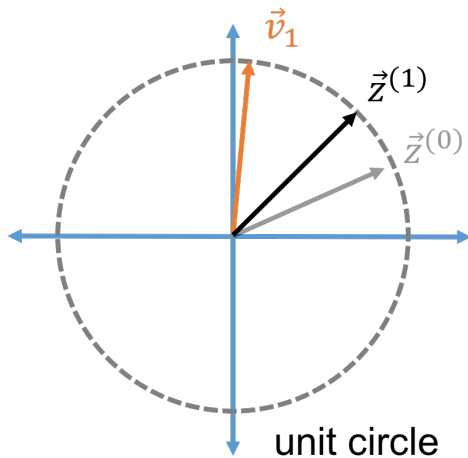
- **Initialize:** Choose $\vec{z}^{(0)}$ randomly. E.g. $\vec{z}^{(0)}(i) \sim \mathcal{N}(0, 1)$.
- For $i = 1, \dots, t$
 - $\vec{z}^{(i)} := \mathbf{A} \cdot \vec{z}^{(i-1)}$
 - $\vec{z}_i := \frac{\vec{z}^{(i)}}{\|\vec{z}^{(i)}\|_2}$

Return \vec{z}_t

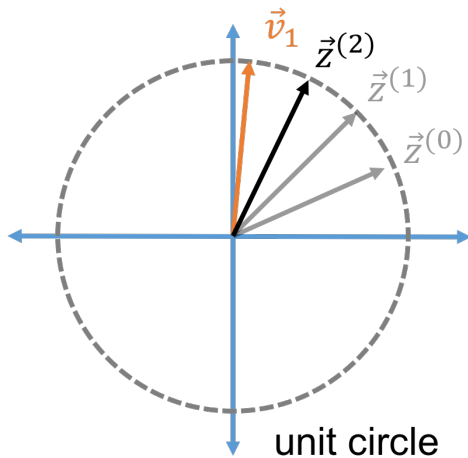
POWER METHOD



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Write $\vec{z}^{(0)}$ in \mathbf{A} 's eigenvector basis:

$$\vec{z}^{(0)} = c_1 \vec{v}_1 + c_2 \vec{v}_2 + \dots + c_d \vec{v}_d.$$

$\mathbf{A} \in \mathbb{R}^{d \times d}$: input matrix with eigendecomposition $\mathbf{A} = \mathbf{V}\mathbf{\Lambda}\mathbf{V}^T$. \vec{v}_1 : top eigenvector, being computed, $\vec{z}^{(i)}$: iterate at step i , converging to \vec{v}_1 .

POWER METHOD ANALYSIS

Write $\vec{z}^{(0)}$ in \mathbf{A} 's eigenvector basis:

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Update step: $\vec{z}^{(i)} = \mathbf{A} \cdot \vec{z}^{(i-1)} = \mathbf{V} \mathbf{\Lambda} \mathbf{V}^T \cdot \vec{z}^{(i-1)}$ (then normalize)

$$\mathbf{V}^T \vec{z}^{(0)} =$$

$$\mathbf{\Lambda} \mathbf{V}^T \vec{z}^{(0)} =$$

$$\vec{z}^{(1)} = \mathbf{V} \mathbf{\Lambda} \mathbf{V}^T \cdot \vec{z}^{(0)} =$$

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Claim 1 : Writing $\vec{z}^{(0)} = c_1 \vec{v}_1 + c_2 \vec{v}_2 + \dots + c_d \vec{v}_d$,

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POWER METHOD CONVERGENCE

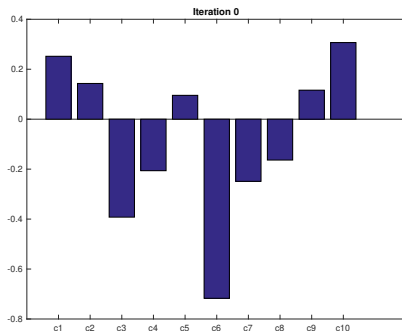
After t iterations, we have 'powered' up the eigenvalues, making the component in the direction of v_1 much larger, relative to the other components.

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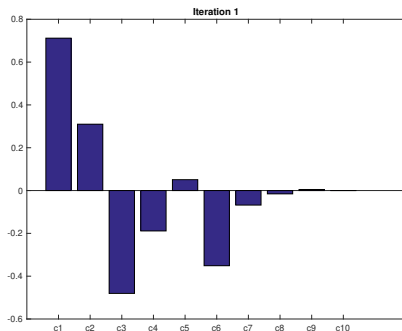
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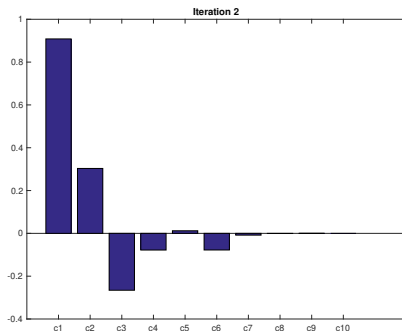
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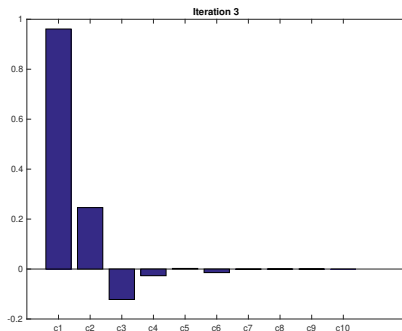
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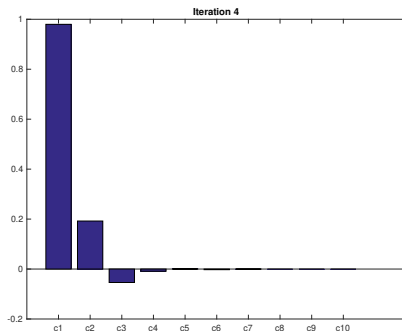
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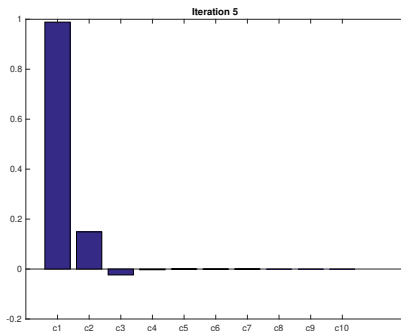
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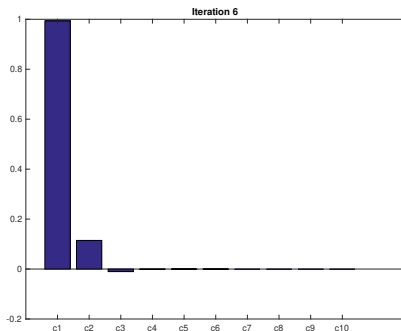
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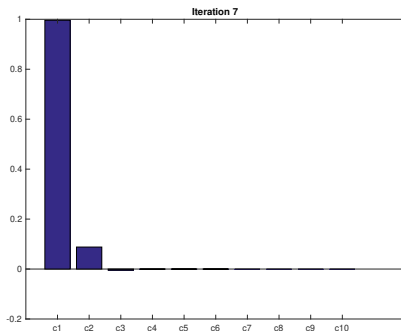
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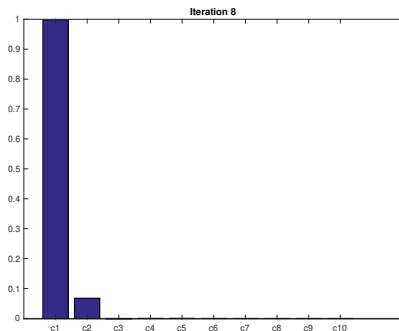
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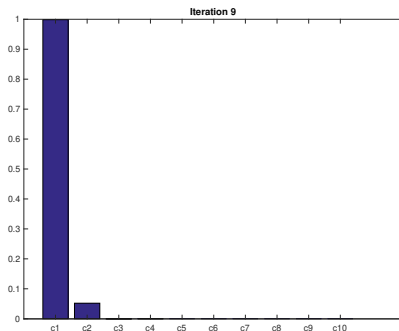
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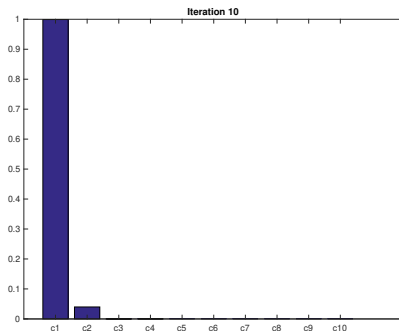
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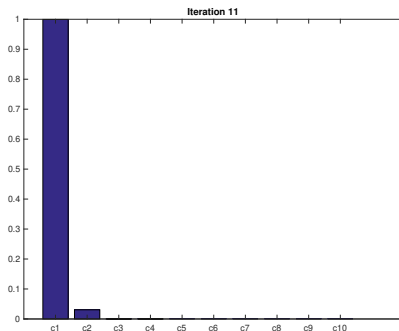
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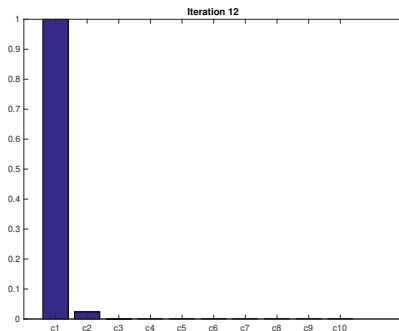
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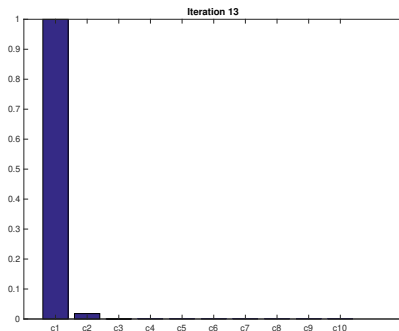
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When will convergence be slow?

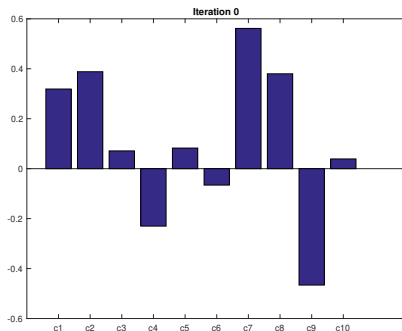
Slow Case: \mathbf{A} has eigenvalues: $\lambda_1 = 1, \lambda_2 = .99, \lambda_3 = .9, \lambda_4 = .8, \dots$

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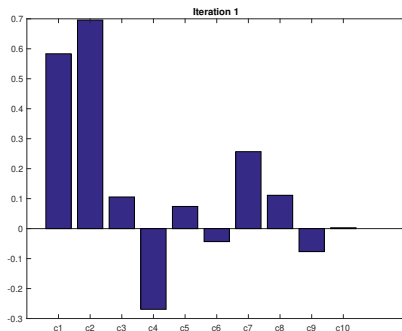
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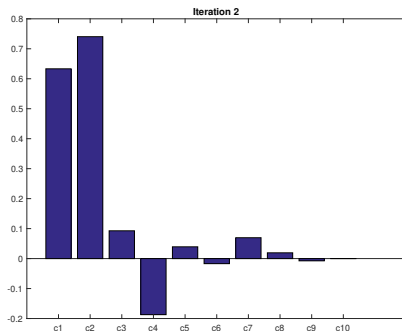
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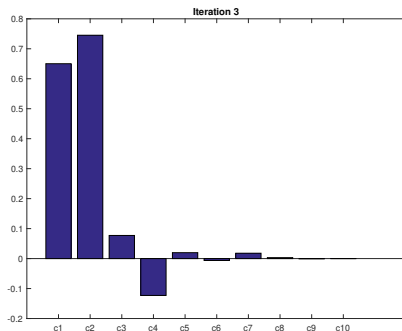
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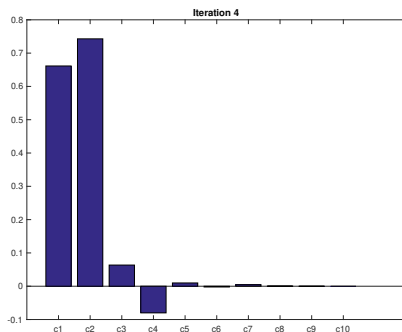
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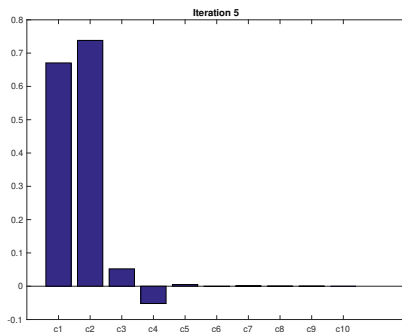
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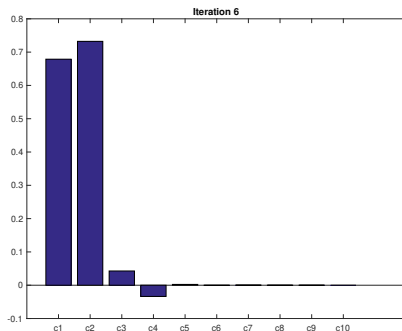
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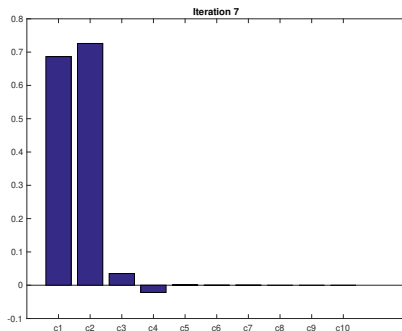
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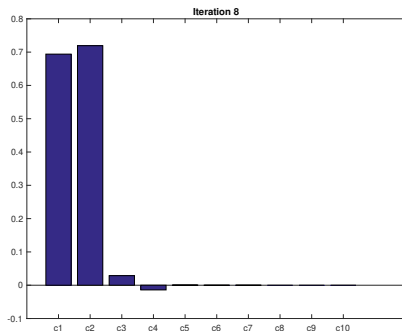
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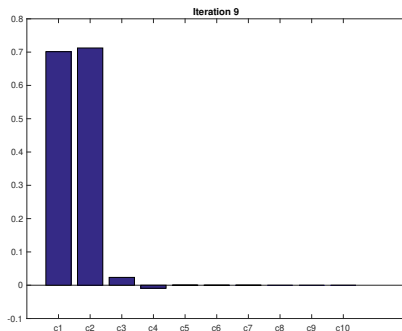
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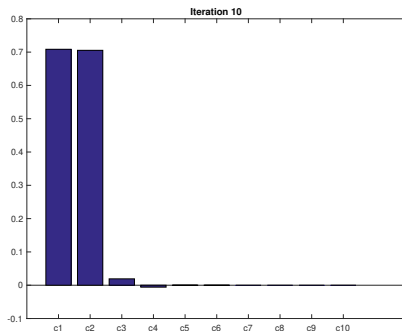
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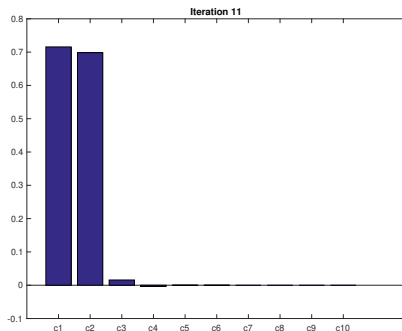
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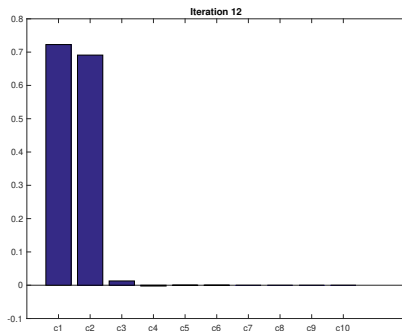
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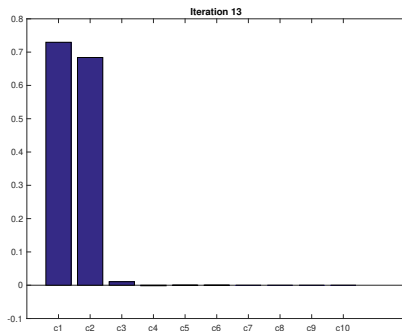
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Write $|\lambda_2| = (1 - \gamma)|\lambda_1|$ for 'gap' $\gamma = \frac{|\lambda_1| - |\lambda_2|}{|\lambda_1|}$.

How many iterations t does it take to have $|\lambda_2|^t \leq \frac{1}{e} \cdot |\lambda_1|^t$?

$\mathbf{A} \in \mathbb{R}^{d \times d}$: input matrix with eigendecomposition $\mathbf{A} = \mathbf{V}\mathbf{\Lambda}\mathbf{V}^T$. \vec{v}_1 : top eigenvector, being computed, $\vec{z}^{(i)}$: iterate at step i , converging to \vec{v}_1 .

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How small must we set δ to ensure that $c_1 \lambda_1^t$ dominates all other components and so $\vec{z}^{(t)}$ is very close to \vec{v}_1 ?

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Claim: When $z^{(0)}$ is chosen with random Gaussian entries, writing $z^{(0)} = c_1 \vec{v}_1 + c_2 \vec{v}_2 + \dots + c_d \vec{v}_d$, with very high probability, for all i :

$$O(1/d^2) \leq |c_i| \leq O(\log d)$$

Corollary:

$$\max_j \left| \frac{c_j}{c_1} \right| \leq O(d^2 \log d).$$

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Setting $\delta = O\left(\frac{\epsilon}{d^3 \log d}\right)$ gives $\|\bar{z}^{(t)} - \vec{v}_1\|_2 \leq \epsilon$.

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Theorem (Basic Power Method Convergence)

Let $\gamma = \frac{|\lambda_1| - |\lambda_2|}{|\lambda_1|}$ be the relative gap between the first and second eigenvalues. If Power Method is initialized with a random Gaussian vector $\vec{v}^{(0)}$ then, with high probability, after $t = O\left(\frac{\ln(d/\epsilon)}{\gamma}\right)$ steps:

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POWER METHOD THEOREM

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Can actually do it in $O\left(\frac{\ln(d/\epsilon)}{\sqrt{\gamma}}\right)$ iterations using “Krylov subspace methods” such as Lanczos method and Arnoldi method.

FINDING SECOND (ETC.) EIGENVECTOR

- If A has eigenvectors v_1, \dots, v_n with eigenvalues $\lambda_1, \dots, \lambda_n$ ($|\lambda_1| \geq \dots \geq |\lambda_n|$) then

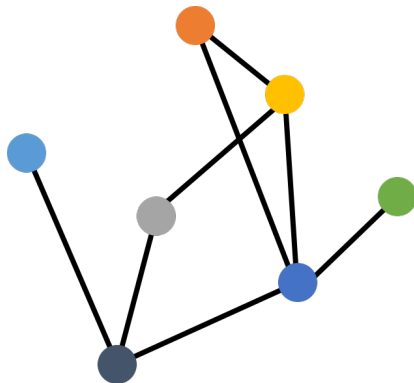
$$B = A - \lambda_1 v_1 v_1^T$$

has eigenvectors v_2, \dots, v_n, v_1 with eigenvalues $\lambda_2, \dots, \lambda_n, 0$

- Hence, to find the second eigenvector of A , just apply the previous method to B .

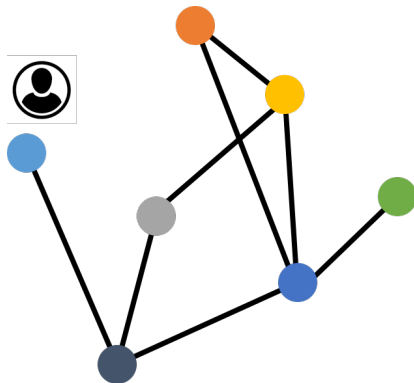
CONNECTION TO RANDOM WALKS

Consider a random walk on a graph G with adjacency matrix \mathbf{A} .



CONNECTION TO RANDOM WALKS

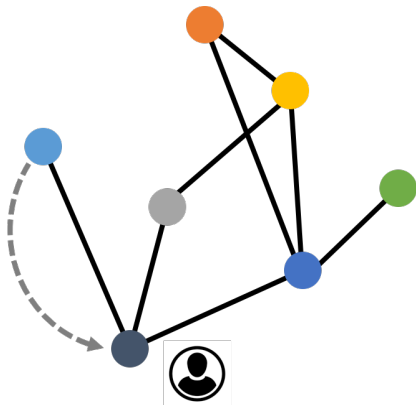
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At each step, move to a random vertex, chosen uniformly at random from the neighbors of the current vertex.

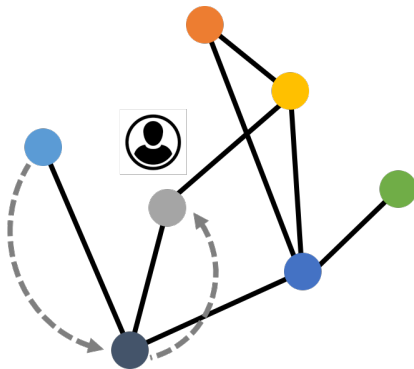
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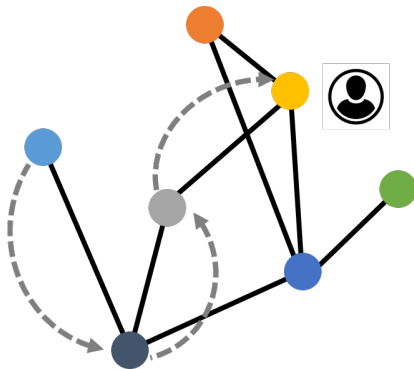
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- **Update:**

$$\Pr(\text{walk at } i \text{ at step } t) = \sum_{j \in \text{neigh}(i)} \Pr(\text{walk at } j \text{ at step } t-1) \cdot \frac{1}{\text{degree}(j)}$$

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Claim: After t steps, the probability that a random walk is at node i is given by the i^{th} entry of

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- Like the power method, the time a random walk takes to converge to its stationary distribution (mixing time) is dependent on the gap between the top two eigenvalues of $\mathbf{D}^{-1/2}\mathbf{A}\mathbf{D}^{-1/2}$. The **spectral gap**.