THE UNIVERSITY OF TEXAS AT AUSTIN
Department of Statistics and Data Sciences
College of Natural Sciences

SDS 385: Stat Models for Big Data

Lecture 10: Pagerank and related methods

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Ranking and Pagerank

- Goal: obtain a ranking of webpages which are connected via hyperlinks
- Hope: webpages pointed to by other "important" webpages are also important.
- Developed by Brin and Page (1999)
- Many subsequent works:
  - HITS (Kleinberg, 1998)
  - Pagerank (Page and Brin, 1998)

Definitions

- $n \times n$ Adjacency matrix A
  - A_{ij} = weight on an edge from i to j
  - If graph is undirected $A(i,j) = A(j,i)$
- $n \times n$ Probability transition matrix P
  - P has rows summing to one, i.e. row stochastic
  - $P(i,j)$ is the probability that a random walker will step on j from i.
  - $P(i,j) = \dfrac{A(i,j)}{\sum_j A(i,j)}$
- $n \times n$ Laplacian matrix L
  - $L = D - A$, where D is the diagonal matrix of degrees
  - It is symmetric positive semidefinite for undirected graphs.
  - Singular, i.e. has a zero eigenvalue

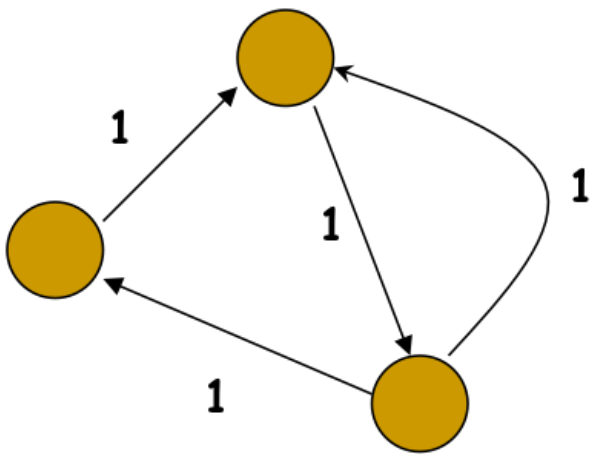
Definition

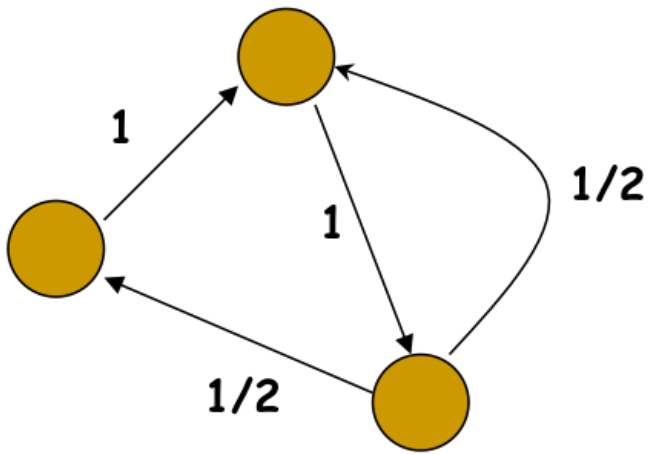
0	1	0
0	0	1
1	1	0

Adjacency matrix A

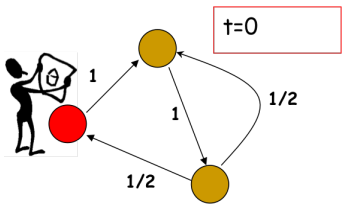
0	1	0
0	0	1
1/2	1/2	0

Transition matrix P





Random walks



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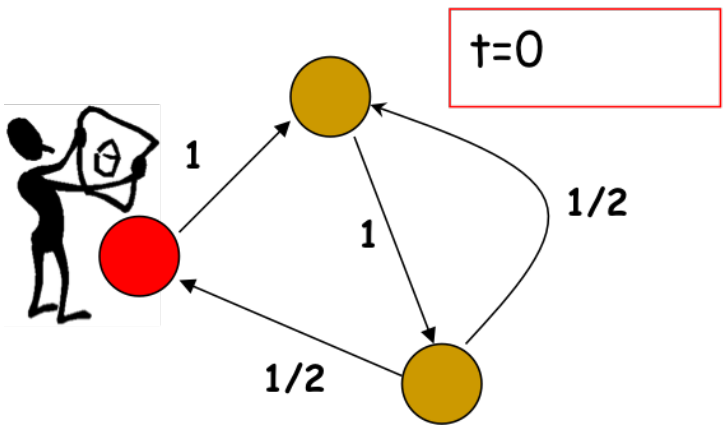

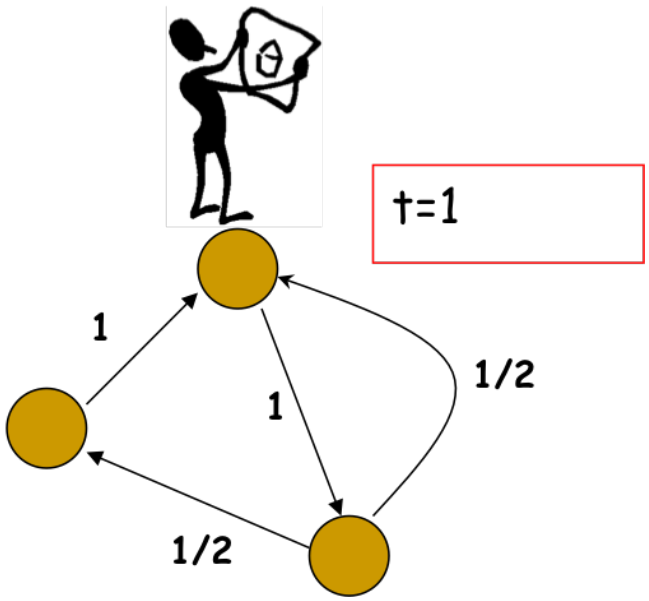

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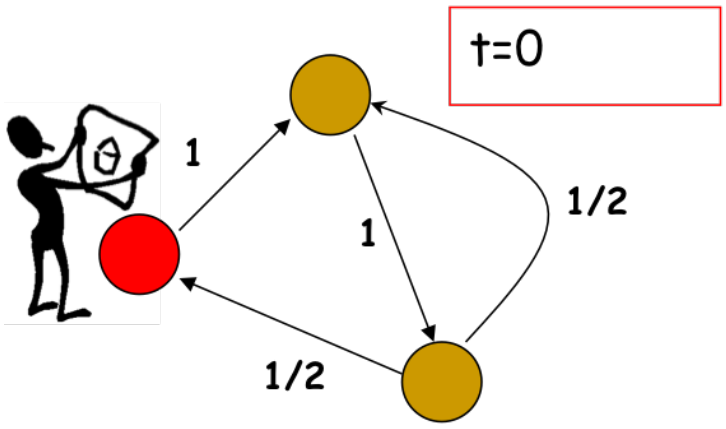

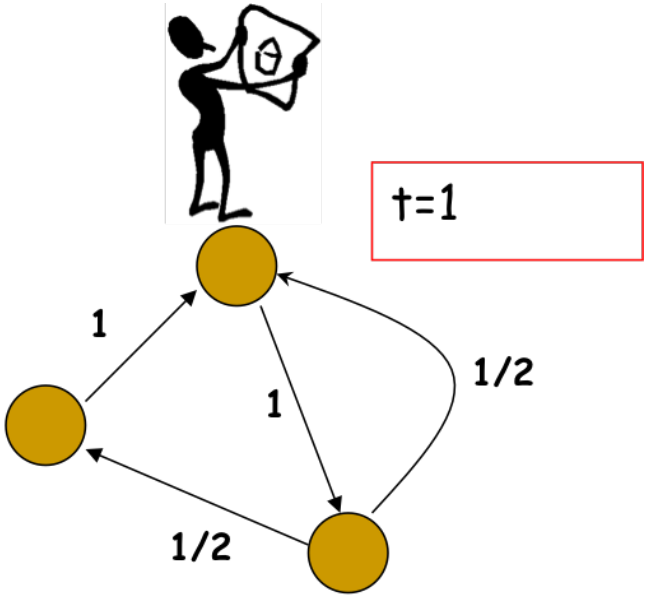

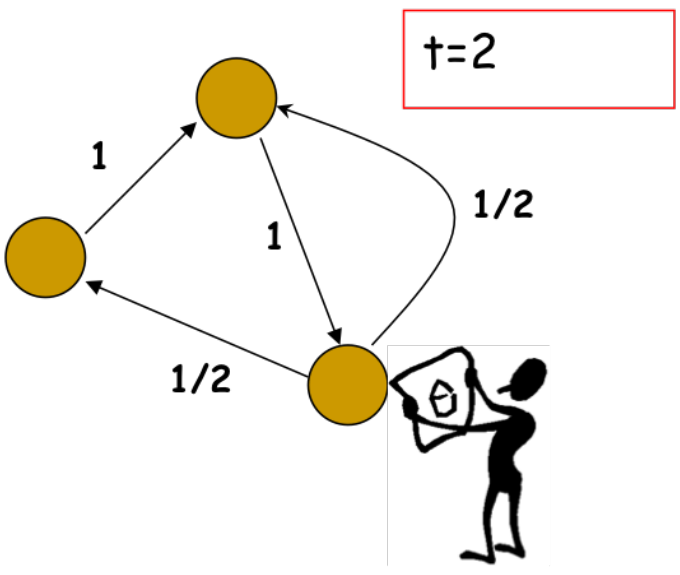

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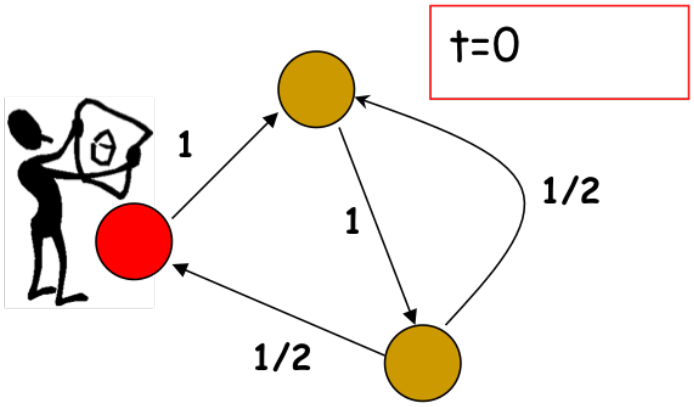

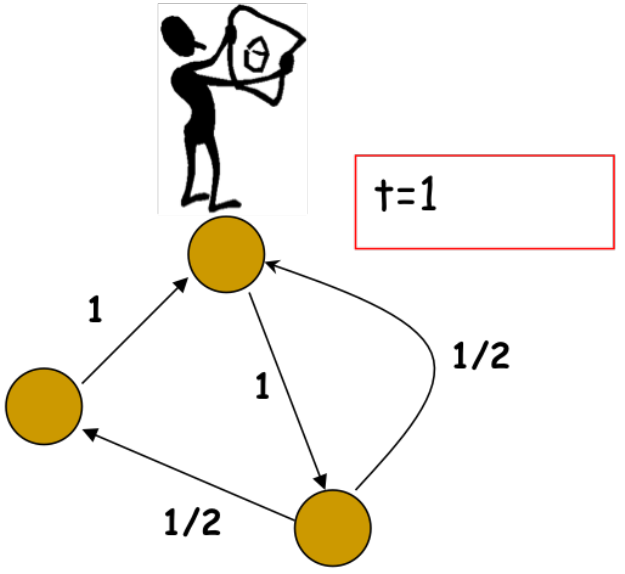

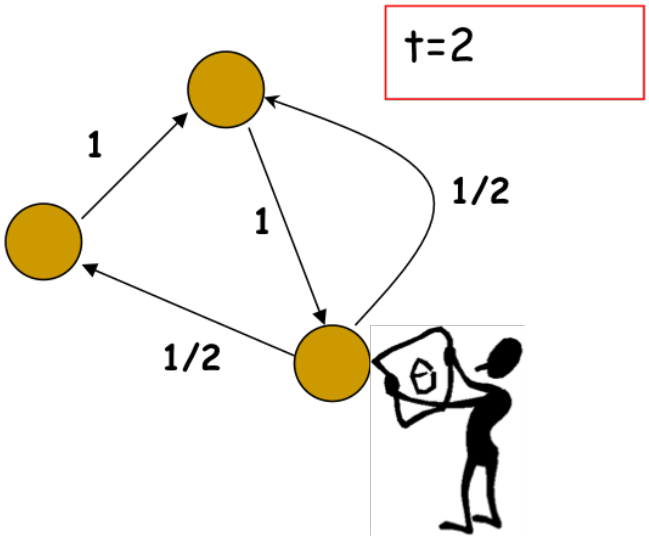

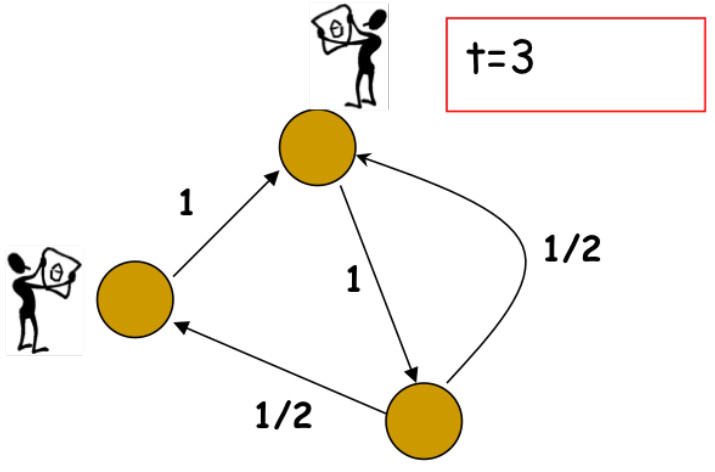

Probability Distributions

- $x_t(i)$ denotes the probability that the surfer is at node i at time t.

$$x_{t+1}(i) = \sum_j x_t(j) P(j, i)$$

$$x_{t+1}^T = x_t^T P = x_{t-1}^T P^2 = \cdots = x_0^T P^t$$

- What happens if the surfer keeps walking for a long time?

Stationary Distribution

- When the surfer keeps walking for a long time
- When the distribution does not change anymore i.e. $x_{T+1} = x_T$
- For well-behaved graphs this does not depend on the start distribution!!

Stationary distribution

- The stationary distribution at a node is related to the amount of time a random walker spends visiting that node.

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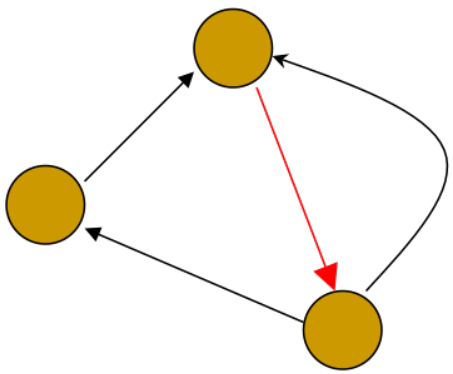
- Whoa! thats just the left eigenvectorof the transition matrix !

Stationary distribution

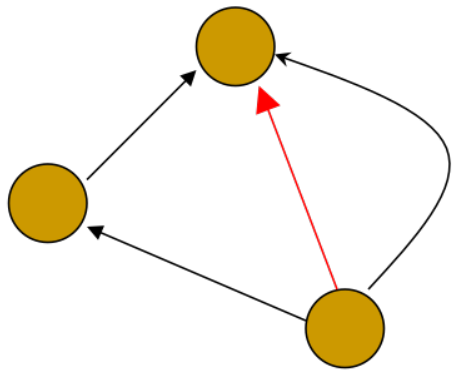
- Lot of theory hiding here.
- For example, what is the guarantee that there will be a unique left eigenvector, or the random walk will at all converge?
- Can't it just keep oscillating?

Well-behaved Markov chains

- Irreducible: There is a path from every node to every other node.



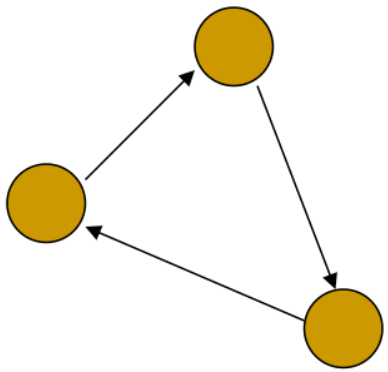
Irreducible



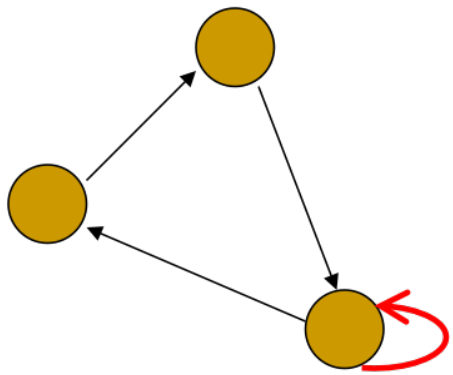
Not irreducible

Well-behaved Markov chains

- Aperiodic: The GCD of all cycle lengths is *1*. The GCD is also called period.



Periodicity is 3



Aperiodic

Well-behaved Markov chains

- If a markov chain is irreducible and aperiodic then the largest eigenvalue of the transition matrix will be equal to 1 and all the other eigenvalues will be strictly less than 1.
- These results imply that for a well behaved graph there exists an unique stationary distribution.

Pagerank and Perron Frobenius

- Perron Frobenius only holds if the graph is irreducible and aperiodic.
- But how can we guarantee that for the web graph? Do it with a small restart probability c.
- At any time-step the random surfer
  - jumps (teleport) to any other node with probability c
  - jumps to its direct neighbors with total probability $1 - c$.

$$\tilde{P} = (1 - c)P + c11^T/n$$

Power iteration

- Power Iteration is an algorithm for computing the stationary distribution.
  - Start with any distribution x_0
  - Keep computing $x_{t+1}^T = x_t^T P$
  - Stop when x_{t+1} and x_t are almost the same

Power Iteration

- Why should this work?
- Write x_0 as a linear combination of the left eigenvectors $\{v_0, v_1, \ldots, v_{n-1}\}$ of P
- Remember that v_0 is the stationary distribution.

$$x_0 = c_0 v_0 + c_1 v_1 + c_2 v_2 + \ + c_{n-1} v_{n-1}$$

Power Iteration

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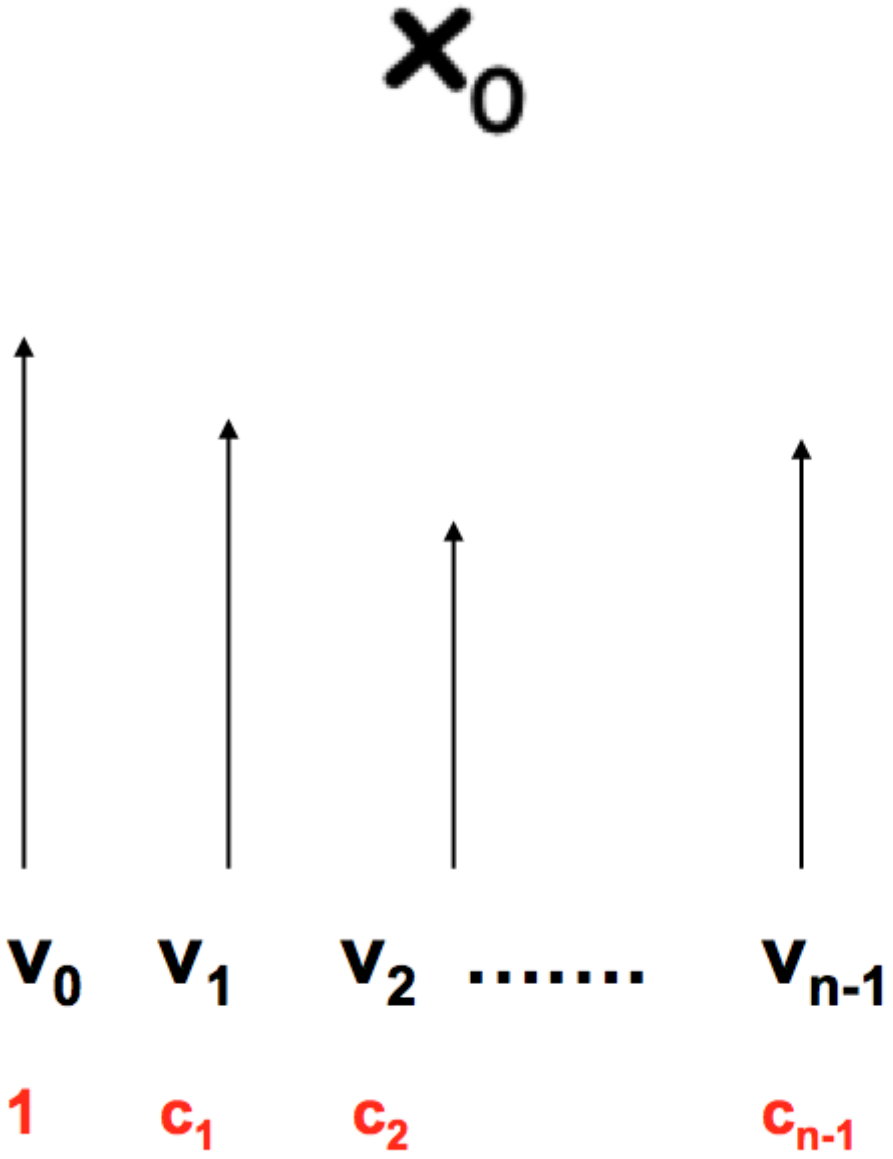
$$x_0 = c_0 v_0 + c_1 v_1 + c_2 v_2 + \ + c_{n-1} v_{n-1}$$

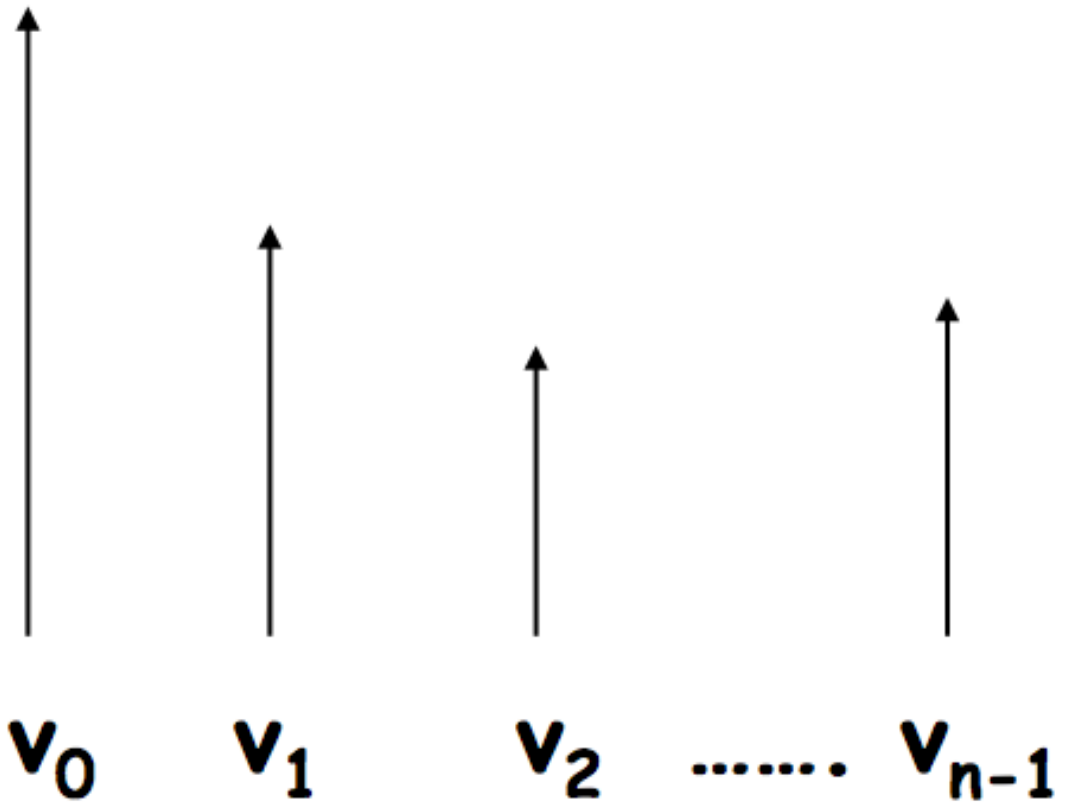
- $c_0 = 1$. Why?
  - First note that $1^T v_i = 0$ if $i \neq 1$
  - So $x_0^T 1 = c_0 = 1$, since both x_0 and v_0 are distributions.

Power Iteration



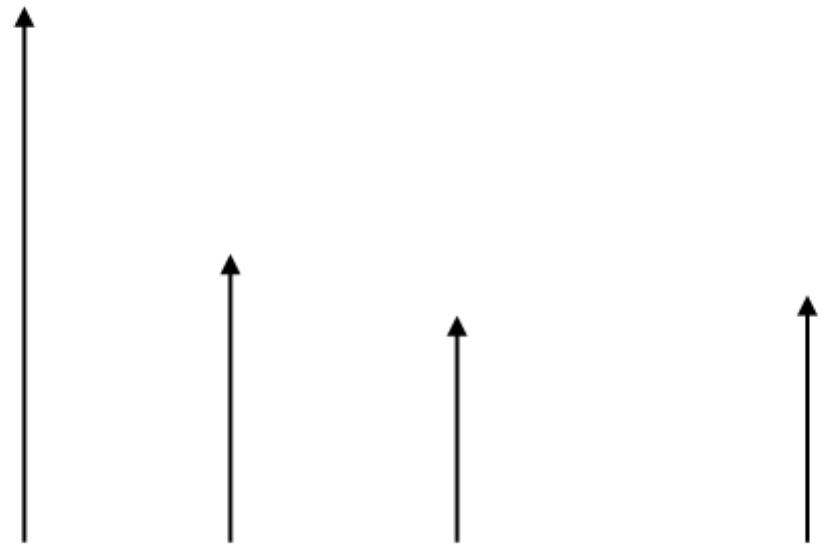
Power Iteration

$$x_1 = x_0 \tilde{P}$$



Power Iteration

$$x_2 = x_1 \tilde{P} = x_0 \tilde{P}^2$$

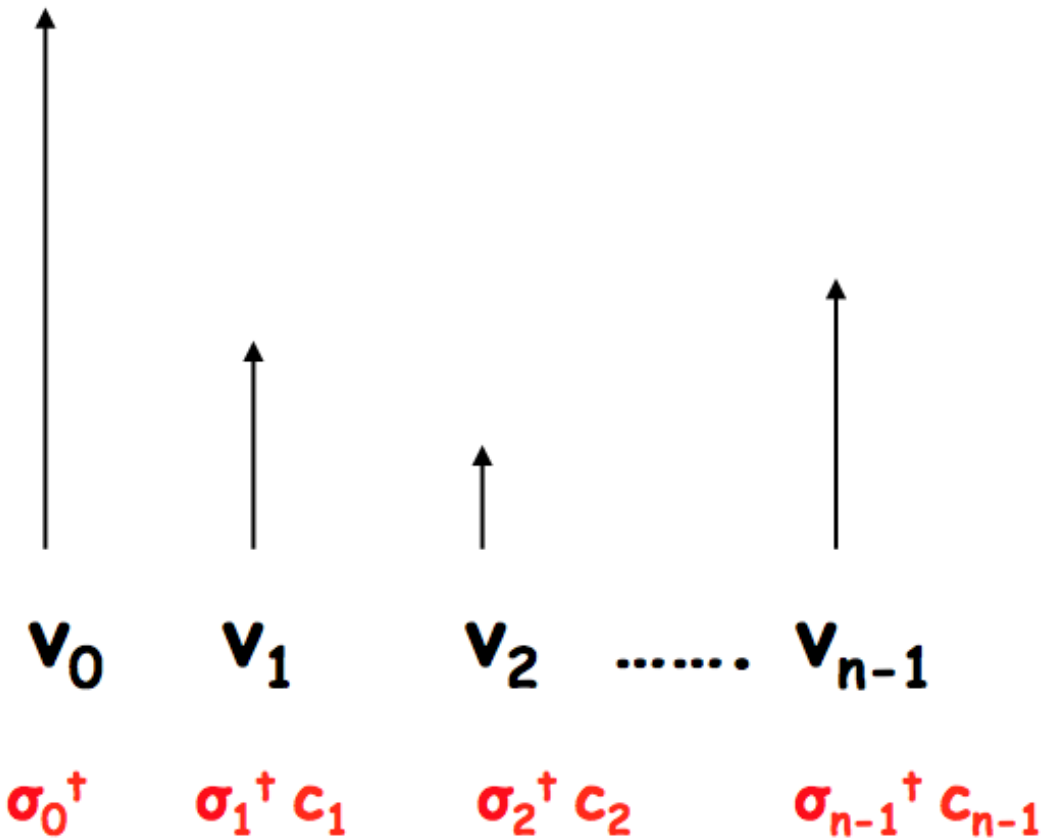


v_0 v_1 v_2 ..... v_{n-1}

σ_0^2 $\sigma_1^2 c_1$ $\sigma_2^2 c_2$ $\sigma_{n-1}^2 c_{n-1}$

Power Iteration

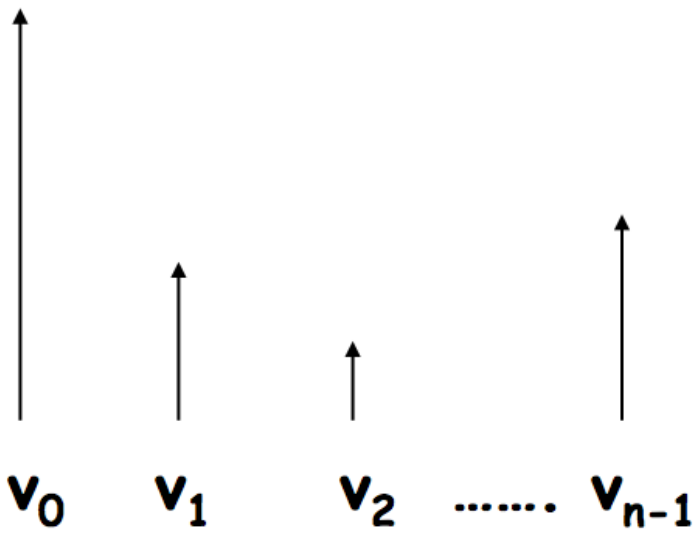
$$x_t = x_0 \tilde{P}^t$$



Power Iteration

$$x_t = x_0 \tilde{P}^t$$

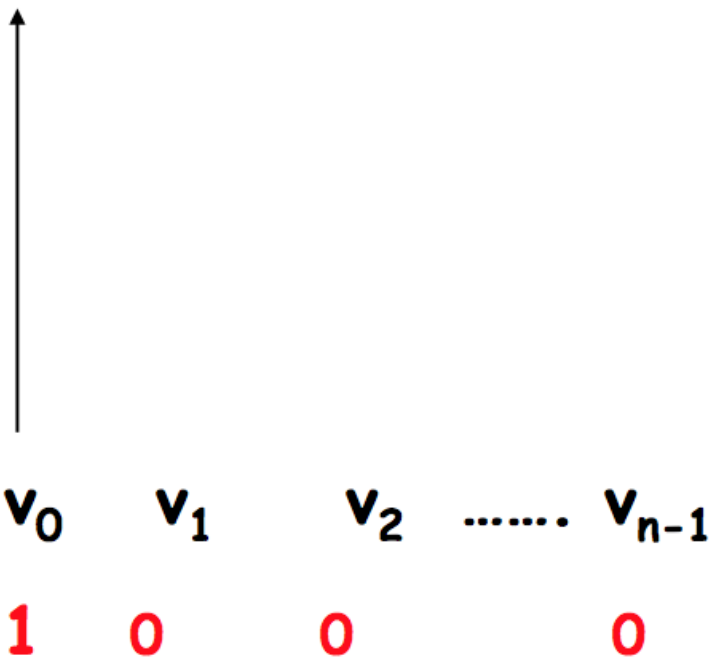
$\sigma_0 = 1 > \sigma_1 \geq \ldots \geq \sigma_n$



Power Iteration

X_∞

$\sigma_0 = 1 > \sigma_1 \geq \ldots \geq \sigma_n$



Second eigenvalue

- Smaller σ_2 is faster the chain mixes.
- For pagerank, we wonder what the second largest eigenvalue is of $\tilde{P} = (1 - c)P + cU$
- The largest eigenvalue is 1
- The second largest is less than $1 - c$ in magnitude.
- So pagerank computation converges fast.

Pagerank

- We are looking for the vector v s.t.

$$v^T = (1-c)v^T P + cr^T$$

- r is a distribution over web-pages.
- If r is the uniform distribution we get pagerank.
- What happens if r is non-uniform?

Pagerank

- We are looking for the vector v s.t.

$$v^T = (1-c)v^T P + cr^T$$

- r is a distribution over web-pages.
- If r is the uniform distribution we get pagerank.
- What happens if r is non-uniform?
- Personalization

Personalized Pagerank

- The only difference is that we use a non-uniform teleportation distribution, i.e. at any time step teleport to a set of webpages.
- In other words we are looking for the vector v s.t.

$$v^T = (1 - c)v^T P + cr^T$$

- r is a non-uniform preference vector specific to an user.
- v gives personalized views of the web.

Personalized Pagerank

- Pre-computation: r is not known from before
- Computing during query time takes too long
- A crucial observation1 is that the personalized pagerank vector is linear w.r.t r

$$\mathbf{r} = \begin{pmatrix} \alpha \\ 0 \\ 1-\alpha \end{pmatrix} \Rightarrow v(r) = \alpha v(r_0) + (1-\alpha) v(r_2)$$

$$\mathbf{r}_0 = \begin{pmatrix} 1 \\ 0 \\ 0 \end{pmatrix}, \mathbf{r}_2 = \begin{pmatrix} 0 \\ 0 \\ 1 \end{pmatrix}$$

- Lots of literature for computing personalized pagerank fast, and on the go.

Rank Stability

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Rank Stability

- How does the ranking change when the link structure changes?
- The web-graph is changing continuously.
- How does that affect page-rank?

Rank Stability

Rank on the entire database.

Rank on 5 perturbed datasets by deleting 30% of the papers

1	"Genetic Algorithms in Search, Optimization and...", Goldberg	1	1	1	1	1
2	"Learning internal representations by error...", Rumelhart+al	2	2	2	2	2
3	"Adaptation in Natural and Artificial Systems", Holland	3	5	6	4	5
4	"Classification and Regression Trees", Breiman+al	4	3	5	5	4
5	"Probabilistic Reasoning in Intelligent Systems", Pearl	5	6	3	6	3
6	"Genetic Programming: On the Programming of ...", Koza	6	4	4	3	6
7	"Learning to Predict by the Methods of Temporal ...", Sutton	7	7	7	7	7
8	"Pattern classification and scene analysis", Duda+Hart	8	8	8	8	9
9	"Maximum likelihood from incomplete data via...", Dempster+al	10	9	9	11	8
10	"UCI repository of machine learning databases", Murphy+Aha	9	11	10	9	10
11	"Parallel Distributed Processing", Rumelhart+McClelland	-	-	-	10	-
12	"Introduction to the Theory of Neural Computation", Hertz+al	-	10	-	-	-

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2. **Automating the contruction of Internet portals with machine learning,** A. Mc Callum, K. Nigam, J. Rennie, K. Seymore, In Information Retrieval Journel, 2000

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Rank Stability

- Ng et al 2001: $\tilde{P} = (1-c)P + cU$

- Theorem: if v is the left eigenvector of . Let the pages $i_1, i_2, \ldots, i_k$ be changed in any way, and let v' be the new pagerank. Then

$$\|v - v'\|_1 \leq \frac{\sum_{j=1}^{k} v(i_j)}{c}$$

Rank Stability

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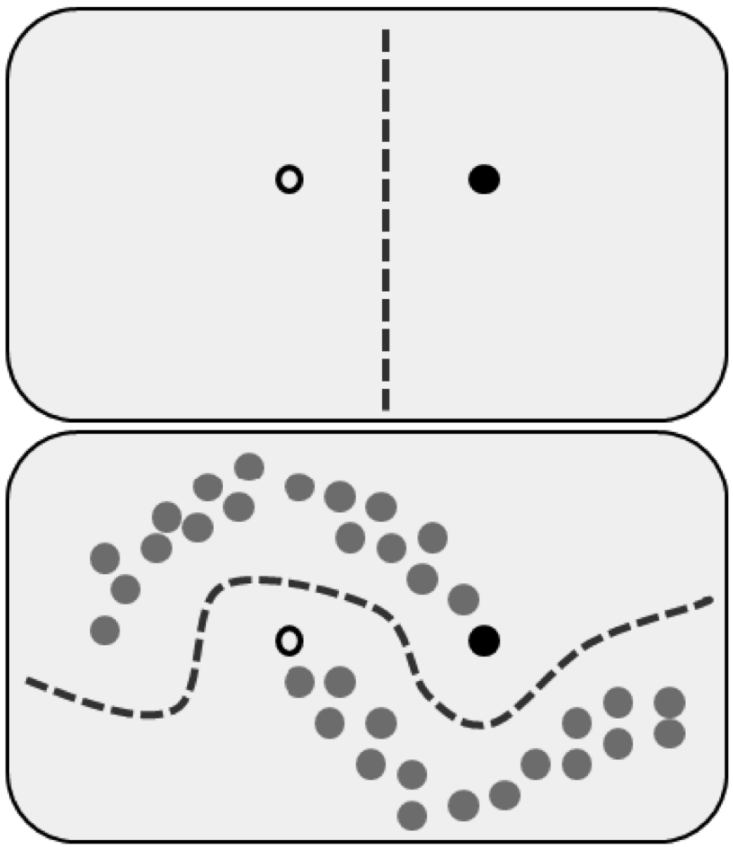
$$\|v - v'\|_1 \leq \frac{\sum_{j=1}^{k} v(i_j)}{c}$$

- So if c is not too close to 0, the system would be rank stable and also converge fast!

Semi-supervised learning

- You are given a lot of unlabeled data.
- Only a few points are labeled.
- Is this useful?



Semi-supervised learning

- Two broad ways
  - Label propagation:
    - Graph Based algorithm
    - Does not generalize to unseen data, i.e. Transductive
  - Manifold regularization
    - Graph Based regularization
    - Does generalize to unseen data, i.e. Inductive

Semi-supervised learning

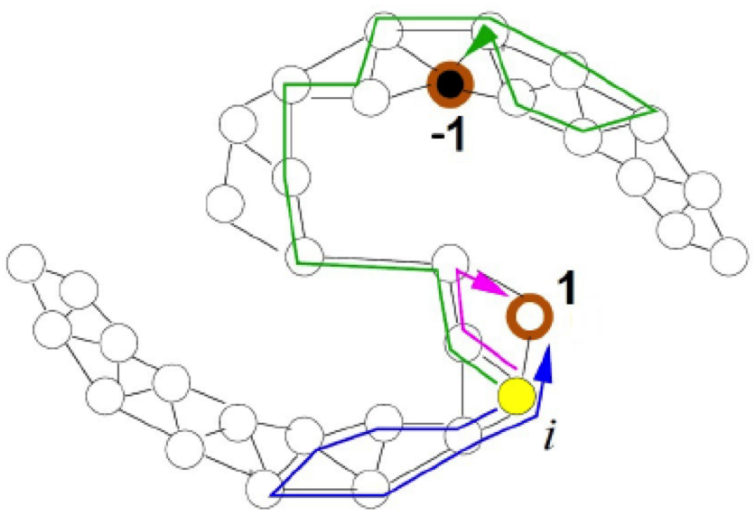
- Input n data points $x_1, \ldots, x_n$
- Define similarity matrix $S \in \mathbb{R}^{n \times n}$

$$S_{ij} = \exp(-\|x_i - x_j\|^2 / 2\sigma^2)$$

- Since S is dense, often k nearest neighbor graphs are also used. (Your homework!)

Label propagation [Zhu et al, 2003]

- Input:
  - ℓ labeled datapoints $(x_1, y_1), \ldots, (x_\ell, y_\ell)$
  - u unlabeled datapoints $x_{\ell+1}, \ldots, x_{\ell+u}$
- Predict: labels $y_{\ell+1}, \ldots, y_{\ell+u}$



Label propagation algorithm

- Compute $P \in [0,1]^{(\ell+u)\times(\ell+u)}$
- $P_{ij} = \dfrac{S_{ij}}{\sum_j S_{ij}}$
- Harmonic function:
  - Function value at an unlabeled node is an average of function values at its neighbors
  - For $j = \ell + 1 : \ell + u$,

$$f(j) = \frac{\sum_i S_{ij} f(i)}{\sum_i S_{ij}}$$

  - In other words, for the unlabeled nodes, this fixed point equation is satisfied

$$f[U] = Pf[U] \qquad f[L] = y[L]$$

  - For a vector v and set S, we denote by $v[S]$ the subset of values in S
  - U and L denote the set of unlabeled and labeled points respectively.

Closed form

- Assume there are just two classes. Set $y[L] \in \{0,1\}^{\ell}$ accordingly.
- We have:

$$\begin{bmatrix} P_{LL} & P_{LU} \\ P_{UL} & P_{UU} \end{bmatrix} \begin{bmatrix} Y_L \\ Y_U \end{bmatrix} = \begin{bmatrix} Y_L \\ Y_U \end{bmatrix}$$

- Expanding, we get:

$$P_{UL} Y_L + P_{UU} Y_U = Y_U$$

- Moving things around:

$$Y_U = (I - P_{UU})^{-1} P_{UL} Y_L$$

- Can use a linear system solver

Label propagation - Random walk interpretation

- Think of the labeled nodes as absorbing states
- Use

$$Y_U = \sum_{t=0}^{\infty} P_{UU}^t P_{UL} Y_L$$

$$= \underbrace{P_{UL} Y_L}_{\text{Probability of reaching "1"s in one step}} + \underbrace{P_{UU} P_{UL} Y_L}_{\text{Probability of reaching in two steps}} +$$

$$= \text{Probability of reaching a label "1" in a long random walk}$$

- Why is this useful?
  - If the labels are all reachable, a long walk must hit a "0" or a "1"
  - So if $Y_i > 1/2$, that means from i, its more likely to reach a "1" than a "0"

Graph Laplacian interpretation

- Graph Laplacian: $L = D - S$, where $D_{ii} = \sum_j S_{ij}$ is a diagonal matrix
- L is positive semi-definite (we are assuming $S_{ij} > 0$)
- Why?
  - For any vector v,
$$v^T L v = \sum_{ij} S_{ij}(v_i - v_j)^2$$
  - So this measures how unsmooth v is w.r.t S
  - L is also singular, why?

Label propagation algorithm

- We can also frame label propagation as

$$\arg\min_{v} v^T L v \qquad s.t. v[L] = y_L$$

- Why?

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- Why?
- Write $v = \begin{bmatrix} y_L & v_U \end{bmatrix}$
- Now set the derivative to zero.
  - $Lv = 0$ such that $v[L] = y_L$
  - $$\begin{bmatrix} D_{LL} - S_{LL} & -S_{LU} \\ -S_{UL} & D_{UU} - S_{UU} \end{bmatrix} \begin{bmatrix} y_L \\ v_U \end{bmatrix} = 0$$
  - Solving:

$$-S_{UL} y_L + (D_{UU} - S_{UU}) v_U = 0$$

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  - Solving:

$$-S_{UL} y_L + (D_{UU} - S_{UU}) v_U = 0$$

  - rearranging:

$$v_U = (D_{UU} - S_{UU})^{-1} S_{UL} y_L = (I - P_{UU})^{-1} P_{UL} y_L$$

Experimentally?

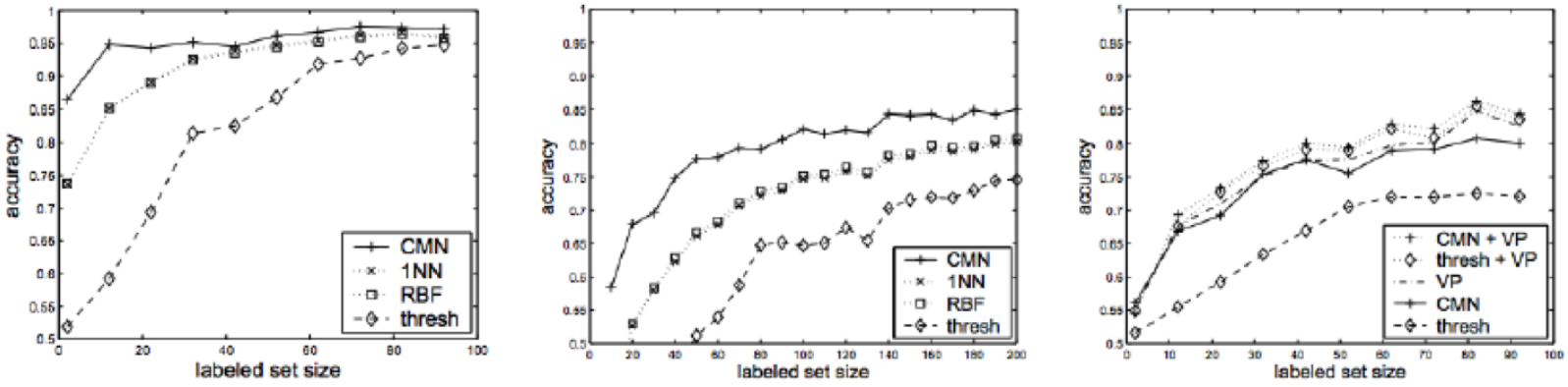

Figure 3. Harmonic energy minimization on digits "1" vs. "2" (left) and on all 10 digits (middle) and combining voted-perceptron with harmonic energy minimization on odd vs. even digits (right)

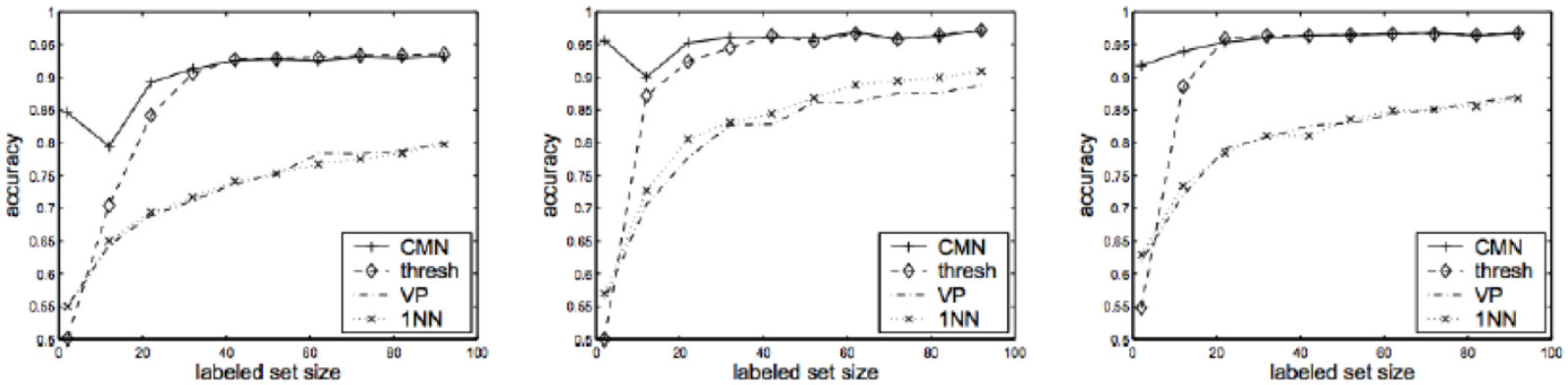

Figure 4. Harmonic energy minimization on PC vs. MAC (left), baseball vs. hockey (middle), and MS-Windows vs. MAC (right)

Manifold regularization

- Input:
  - ℓ labeled datapoints $(x_1, y_1), \ldots, (x_\ell, y_\ell)$
  - u unlabeled datapoints $x_{\ell+1}, \ldots, x_{\ell+u}$
- Predict: labels $y_{\ell+1}, \ldots, y_{\ell+u}$

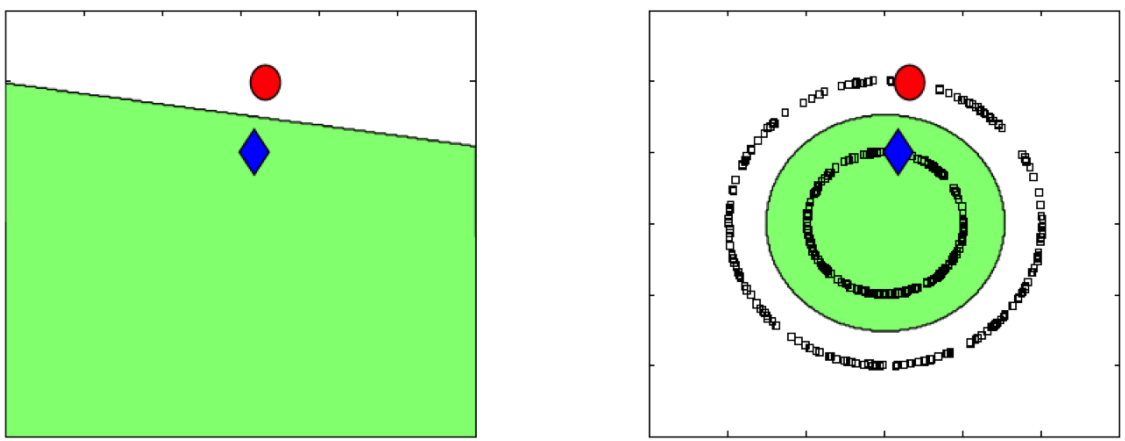

Figure 1: Unlabeled data and prior beliefs

Manifold regularization

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- Before the constraint was $v[L] = y_L$, now instead we will use a loss function and learn a classifier on the labeled data

$$\min_w \underbrace{\sum_{i=1}^{\ell} \text{loss}(y_i, w^T x_i)}_{\textit{loss}} + \lambda \underbrace{R(w)}_{\textit{regularization}}$$

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- How about the unlabeled data?

Manifold regularization: Belkin et al 2006

$$\min_w \underbrace{\sum_{i=1}^{\ell} \text{loss}(y_i, w^T x_i)}_{loss} + \lambda \underbrace{R(w)}_{regularization} + \beta (Xw)^T L(Xw)$$

- Assume a linear predictor $w^T x$
- Idea: close/similar points have similar predicted labels.
- LapSVM:

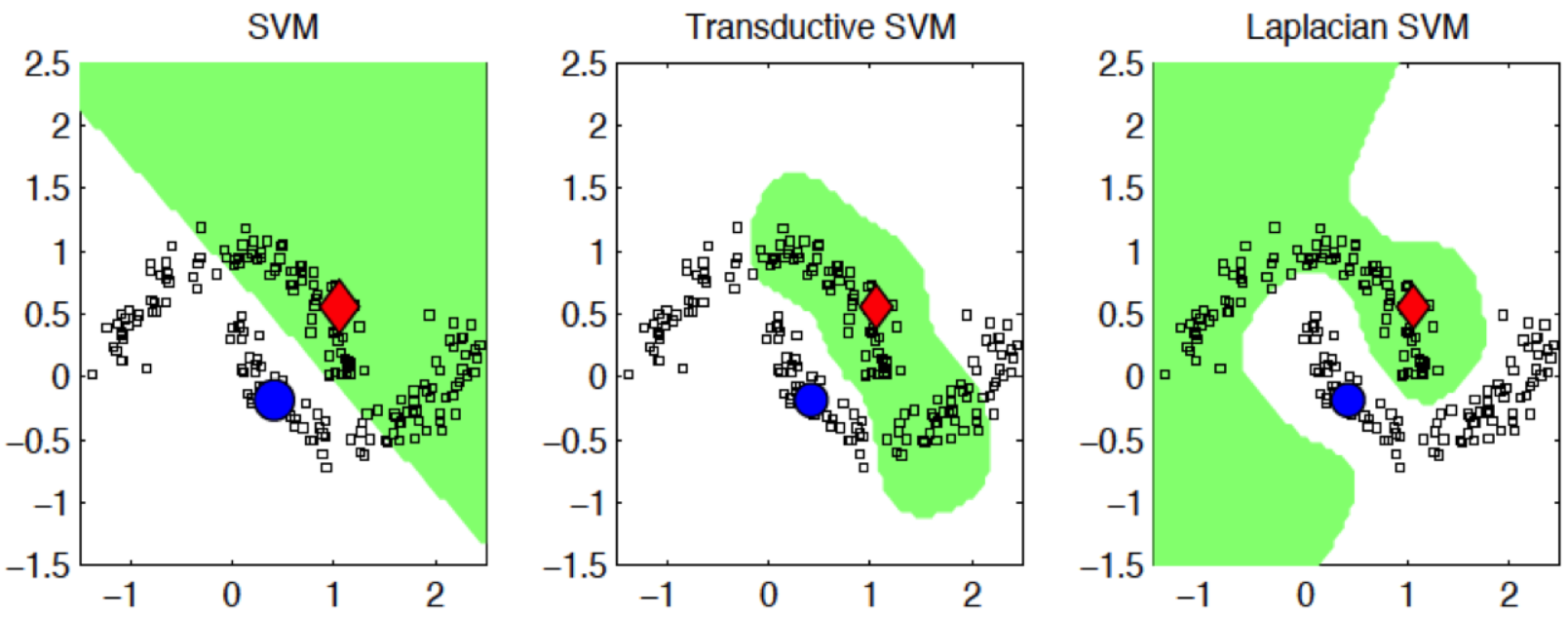
$$\min_w \sum_{i=1}^{\ell} (1 - y_i f(x_i))_+ + \lambda \underbrace{\|f\|_K^2}_{regularization} + \beta f^T L f$$

Transductive SVM: Joachims et al 1999

$$\min_{w, y_{\ell+1}, \ldots, y_{\ell+u}} \sum_{i=1}^{\ell} (1 - y_i f(x_i))_+ + C' \sum_{i=\ell+1}^{u} (1 - y_i f(x_i))_+ \lambda \underbrace{\|f\|_K^2}_{\textit{regularization}}$$

- Iteratively solves SVM quadratic programs
- Switches labels to improve objective function
- Suffers from local optima, inherently combinatorial problem

Transductive SVM VS LapSVM



Summary

- Three main parts$+\epsilon$
- Large scale optimization:
  - Gradient descent, Newton Raphson
  - Stochastic gradient descent, proximal methods, subgradients, dual coordinate ascent, etc.
    - Le Cun, Leon Bottou Yann. "Large Scale Online Learning." NIPS, 2004.
    - Main point: Given the same time of training, SGD can allow more datapoints, and hence may have faster convergence

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    - John Duchi, Elad Hazan, Yoram Singer. "Adaptive Subgradient Methods for Online Learning and Stochastic Optimization." Journal of Machine Learning Research 12 (2011): 2121-2059.
    - Use the geometry of the data to choose a better loss function. Bottomline - use diagonal approximation of the Hessian instead of full Hessian.

Summary

- Large scale optimization:
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    - Rie Johnson and Tong Zhang. Accelerating stochastic gradient descent using predictive variance reduction. In Advances in neural information processing systems, pages 315323, 2013
    - Main point: Talks about dual coordinate ascent and shows how this leads to variance reduction

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    - Yu et al., Large Linear Classication when Data Cannot Fit In Memory", in KDD 2010) Oct 10
    - Block updates by dividing data into chunks that fit into main memory.

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  - Weinberger, Kilian, et al. "Feature hashing for large scale multitask learning." ICML, 2009. Oct 17
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  - Doing Johnson Lindenstrauss with very sparse projection matrices
  - Gray, Alexander, Ting Liu and Andrew W. Moore. "New Algorithms for Efficient High-Dimensional Nonparametric Classification." Journal of Machine Learning Research 7 (2006): 1135-1158.
  - Fast tree based methods for nearest neighbor search when data has imbalanced classes

Summary

- Parallel methods and Divide and Conquer: better compression for parallel SGD. Divide and conquer methods for community detection in very large graphs. Bootstrap and subsampling- Bag of little bootstraps.
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  - Asynchronous SGD without locks–use the sparsity in data
  - Mackey, Lester, Ameet Talwalkar, Michael I. Jordan. "Divide-and-Conquer Matrix Factorization." arXiv:1107.0789v6 (2012)
  - Divide up the data, do matrix factorization on each part and do a

Summary

- The ϵ - Semisupervised learning and Pagerank. Some Bayesian methods.
    - Jasper Snoek, Hugo Larochelle, and Ryan P Adams. "Practical bayesian optimization of machine learning algorithms"