Accelerated and Sparse Algorithms for Personalized PageRank

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joint work with Elias Wirth and Sebastian Pokutta (art by DALL-E 3)

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What PageRank is about

Rank all websites on the public internet.



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But also...



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Rank all websites on the public internet.

But also...

Cluster computation: large intraconnectivity, low connectivity with the rest.

Other data analyses on large graphs.



David F. Gleich



The Random Walker

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- In PageRank websites are sorted by importance.
- Rank ^{prop.} visit frequency of a random surfer on the internet.
- The random surfer moves uniformly at random from one page to the next one following hyperlinks.
- This is almost perfect. But we need to fix one issue.





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irreducible

Fundamental Theorem of Markov Chains:

A finite strongly connected Markov chain has a unique stationary distribution π , where $\pi_j = 1/h_{j,j}$, for each state j.

Perron-Frobenius theorem ; If also aperiodic, then $\lim_{n\to\infty} x_0^\top Q^n = \pi$; Sampling ; MCMC]

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- Its transition matrix is $Q = (1 \alpha)AD^{-1} + \alpha \mathbb{1}s^{\top}$.
- Google originally used uniform s = 1/n and $\alpha \approx 0.15$.



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PageRank Problem:

For Q stochastic and irreducible, compute $\pi \in \Delta^n$ such that $\pi^\top Q = \pi$.

- ▶ It is an eigenvector problem, a Principal Component Analysis (PCA) problem.
- But it has structure.

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- A is symmetric, so x^T(1 − α)AD⁻¹ + αe_i = x for x ∈ Δⁿ can be cast as the quadratic minimization:

$$\min_{x\geq 0}\left\{\frac{1}{2}x^{\top}x-\frac{1}{2}x^{\top}((1-\alpha)AD^{-1})x-\alpha x^{\top}e_i\right\}.$$



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• Adding ℓ_1 regularization makes the solution sparse:

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Massage the problem and you obtain, for symmetric Q s.t. $0 \prec \alpha \cdot I \preccurlyeq Q \preccurlyeq I$ and $Q_{ij} \leq 0$ for $i \neq j$:

$$\min_{\mathsf{x}\in\mathbb{R}^n_{\geq \mathbf{o}}}\{g(\mathsf{x})\stackrel{\text{\tiny def}}{=}\frac{1}{2}\langle Q\mathsf{x},\mathsf{x}\rangle-\langle\mathsf{b},\mathsf{x}\rangle\}.$$

ℓ_1 -regularization induces sparsity



Gradient Descent

• Gradient descent with step-size $\leq 1/\lambda_{max}$ stays in the current eigenvector's orthant.

We can show that GD restricted to the vertical axis is always coordinatewise lower than its minimizer.

We can discover good coordinates one by one!



Coordinate discovery by approximate optimization

- 1. Because $Q_{ij} \leq 0$ for $i \neq j$, for $y = x + \Delta e_i$, we have $\forall j \neq i$:
 - $\nabla_j g(y) \le \nabla_j g(x)$ if $\Delta > 0$ $\nabla_j g(y) \ge \nabla_j g(x)$ otherwise.



Coordinate discovery by approximate optimization

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∇_jg(y) ≤ ∇_jg(x) if Δ > 0
∇_jg(y) ≥ ∇_jg(x) otherwise.
Recall, ∇_ig(x^{(*,C^(t))}) < 0 only if i is good. So by 1., for x ∈ C^(t) s.t. x ≤ x^{(*,C^(t))}, new coordinates i can only satisfy ∇_ig(x) < 0 if they are good.



Figure: A negative coordinate gradient for a point $x \le x^{(*,C^{(t)})}$ implies the coordinate is good, but not necessarily if $x \le x^{(*,C^{(t)})}$

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- 2. Recall, $\nabla_i g(\mathbf{x}^{(*,C^{(t)})}) < 0$ only if *i* is good. So by 1., for $x \in C^{(t)}$ s.t. $x \leq \mathbf{x}^{(*,C^{(t)})}$, new coordinates *i* can only satisfy $\nabla_i g(x) < 0$ if they are good.
- Strategy: Get close to x^{(*,C^(t))} with Proj. AGD and then move slightly towards 0 to be ≤ x^{(*,C^(t))}.



Figure: A negative coordinate gradient for a point $x \le x^{(*,C^{(t)})}$ implies the coordinate is good, but not necessarily if $x \le x^{(*,C^{(t)})}$.

Results and comparison

▶ The Hessian of g is Q, satisfying $\mu I \preccurlyeq Q \preccurlyeq LI$, its condition number is L/μ .

► $S^* \stackrel{\text{def}}{=} \operatorname{supp}(x^*)$, $\operatorname{vol}(S^*) \stackrel{\text{def}}{=} \operatorname{nnz}(Q_{:,S^*})$ and $\operatorname{vol}(S^*) \stackrel{\text{def}}{=} \operatorname{nnz}(Q_{S^*,S^*})$.

▶ For ℓ_1 -regularized personalized PageRank, it is vol(δ^*) $\leq \frac{1}{a} + |\delta^*|$ [FRS+19].

Method	Time complexity	Space complexity	$-vol(S^*) \stackrel{\text{def}}{=}$	ş.
ISTA [FRS+19]	$\widetilde{\mathbb{O}}(vol(\mathbb{S}^*)\frac{L}{\mu})$	୦(୫*)	V01(8') =	
CDPR (Ours)	$O(S^* ^3 + S^* vol(S^*))$	𝒴(𝔅* ²)	$\widetilde{vol}(\mathbb{S}^*) \stackrel{\text{\tiny def}}{=}$	(1
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• Time complexity $O(|S^*|^3 + |S^*|vol(S^*))$ and space complexity $O(|S^*|^2)$.

Accelerated and Sparse PageRank (ASPR) algorithm

► Lemma. Let $\bar{\mathbf{x}}^{(t+1)}$ be an $\varepsilon \cdot \frac{\mu^2}{2(1+|\bar{\mathbf{s}}^{(t)}|)L^2}$ minimizer in $C^{(t)}$. Define $\mathbf{x}^{(t+1)} \leftarrow \operatorname{Proj}_{\mathbb{R}^n_{\geq 0}}(\bar{\mathbf{x}}^{(t+1)} - \delta_t \mathbb{1})$ for $\delta_t = \sqrt{\frac{\varepsilon\mu}{(1+|\bar{\mathbf{s}}^{(t)}|)L^2}}$. Then, $\mathbf{x}^{(t+1)} \leq \mathbf{x}^{(*,C^{(t)})}$ and $\mathbf{x}^{(t+1)}$ is a global ε -minimizer or there is i s.t. $\nabla_i g(\mathbf{x}^{(t+1)}) < 0$, so we expand the current set of good coordinates $S^{(t)}$.

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- ▶ Intuition. $x^{(t+1)}$ is almost optimal in $C^{(t)}$, so if its global gap is > ε then 1 step of GD makes more progress than what it is possible in $C^{(t)}$. $\implies \exists i \notin S^{(t)}$ s.t. $\nabla_i g(x^{(t+1)}) < 0$.

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- Subproblem optimization only needs gradients in C^(t), costing O(vol(S*)) each. And one full gradient is used at the end of each stage to find new good coordinates, costing O(vol(S*)). It is done at most |S*| times.

• Time complexity $\widetilde{O}(|S^*|\widetilde{vol}(S^*)\sqrt{\frac{L}{\mu}} + |S^*|vol(S^*))$ and space complexity $O(|S^*|)$.

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CASPR (Ours)	$\widetilde{\mathbb{O}}(\mathbb{S}^* \widetilde{vol}(\mathbb{S}^*)\min\left\{\sqrt{\frac{L}{\mu}}, \mathbb{S}^* \right\}+ \mathbb{S}^* vol(\mathbb{S}^*))$	O(S*)
LASPR (Ours)	$\widetilde{\mathbb{O}}(\mathbb{S}^* vol(\mathbb{S}^*))$	୦(୫ *)

Experiments

Results from a Standford Network Analysis Project graph with 3.7M nodes and 16.5M edges.



Left: Time taken to optimize to 10^{-6} accuracy, while fixing $\rho = 10^{-4}$ and varying the regularization α . Right: Time taken to optimize to 10^{-6} accuracy, while fixing $\alpha = 0.05$ and varying ρ .

Experiments

Results from a Standford Network Analysis Project graph with 3.7M nodes and 16.5M edges.



Left: Gap versus time.

Right: Number of non-zeros of the iterates with time. We obtain greater sparsity. This is due to the algorithms optimizing in the space of currently known good coordinates before adding new ones.

Other research: Packing Proportional Fairness

- Pairs server-user in a shared network with limited link capacities.
- ▶ How much flow should each pair receive, while satisfying proportional fairness axioms?



Other research: Riemannian Optimization in Machine Learning

- ▶ How to approximate $\min_{x \in \mathcal{M}} f(x)$ for a Riemannian manifold \mathcal{M} ?
- Mixture of Gaussians, operator scaling, dictionary learning, low-rank matrix completion, RNNs...



Other research: Decentralized Cooperative Stochastic Bandits

- Decentralized network, each node can only communicate to their neighbors.
- They are facing the same stochastic multi-armed bandit problem. How to behave and share information to minimize regret?





Thank you!

Questions?

