When Can Machines Learn? Insights from Convex Optimization and Semidefinite Programming

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Artificial Intelligence is the New Electricity - Andrew Ng

Machine Learning is the New Alchemy - Ali Rahimi and Ben Recht

Present a semidefinite relaxation for the graph clustering problem based on decomposition of graph into union of disjoint subgraphs.

Give a probabilistic model for **''clusterable''** data and graphs, and theoretical recovery guarantees.

Open problems and current research.

Joint with Polina Bombina (UA) and Aleksis Pirinen (RISE Research Institutes of Sweden).

Supported by NSF Grants #2012554 and #2108645; UA Cyberseed Grant SP14572; University of Alabama RGC grants RG14678 and RG14838. **Clustering**: partition data so that items in each cluster are similar to each other and items not in the same cluster are dissimilar.

Fundamental problem in statistics and machine learning:

• pattern recognition, computational biology, image processing/computer vison, network analysis.

No consensus on what constitutes a **good** clustering; depends heavily on application.

Intractable: usually modeled as some NP-hard problem (e.g., clique, normalized cut, k-means).

Clustering seems to be a very difficult/ill-posed problem.

Many heuristics seem to work well in practice.

Question: can we show that we can cluster "clusterable" data? How do we model clusterable data?

Given data and affinity function f indicating similarity between any two items.

Model the data as weighted similarity graph $G_S = (V, E, W)$ as follows:

- Each item is represented by a node in V.
- We add an edge between each pair of two nodes *i*, *j* with edge weight w_{ij} = f(i, j) ∈ [0, 1].
- *w_{ij}* is large if *i* and *j* are highly similar.

Example: Rehnquist Supreme Court

Data drawn from U.S. Supreme Court decisions (from 1994-95 to 2003-04).

First consider by Hubert and Steinley 2005.

Assign edge-weights corresponding to fraction of decisions on which Justices agreed:

	St	Br	Gi	So	Ос	Ke	Re	Sc	Th
St	1	0.62	0.66	0.63	0.33	0.36	0.25	0.14	0.15
Br	0.62	1	0.72	0.71	0.55	0.47	0.43	0.25	0.24
Gi	0.66	0.72	1	0.78	0.47	0.49	0.43	0.28	0.26
So	0.63	0.71	0.78	1	0.55	0.5	0.44	0.31	0.29
Ос	0.33	0.55	0.47	0.55	1	0.67	0.71	0.54	0.54
Ke	0.36	0.47	0.49	0.5	0.67	1	0.77	0.58	0.59
Re	0.25	0.43	0.43	0.44	0.71	0.77	1	0.66	0.68
Sc	0.14	0.25	0.28	0.31	0.54	0.58	0.66	1	0.79
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We want to partition the graph into cliques with heavy support.

A k-disjoint-clique subgraph of a graph G is a subgraph of G induced by k disjoint cliques.

Densest *k*-disjoint-clique problem (KDC): find a *k*-disjoint-clique subgraph such that the sum of the densities of the *k* complete subgraphs induced by the cliques is maximized.

Density of complete subgraph induced by C:

$$d(C) = \frac{1}{|C|} \sum_{i \in C} \sum_{j \in C} w_{ij} = \frac{\mathbf{v}^T \mathbf{W} \mathbf{v}}{\mathbf{v}^T \mathbf{v}}$$

where \mathbf{v} is the characteristic vector of C.

Let $\{C_1, \ldots, C_k\}$ define a *k*-disjoint-clique subgraph with characteristic vectors $\{\mathbf{v}_1, \mathbf{v}_2, \ldots, \mathbf{v}_k\}$

Lift the *k* characteristic vectors $\{\mathbf{v}_1, \mathbf{v}_2, \dots, \mathbf{v}_k\}$ to the rank-*k* matrix variable **X**:

$$\boldsymbol{X} = \sum_{i=1}^{k} \frac{\boldsymbol{v}_i \boldsymbol{v}_i^T}{\|\boldsymbol{v}_i\|^2} = \sum_{i=1}^{k} \frac{\boldsymbol{v}_i \boldsymbol{v}_i^T}{|C_i|}$$

Want to find **X** that maximizes

$$\operatorname{tr}(\boldsymbol{W}\boldsymbol{X}) = \sum_{i=1}^{k} \frac{\boldsymbol{v}_{i}^{\mathsf{T}} \boldsymbol{W} \boldsymbol{v}_{i}}{\|\boldsymbol{v}_{i}\|^{2}} = \sum_{i=1}^{k} d(C_{i})$$

Lifted solutions

Lifted solution **X** must satisfy:



Inlier rows sum to 1. Outlier rows equal 0: $Xe \leq e$

 $m{X}$ is symmetric doubly nonnegative: $m{X} \ge m{0}, \quad m{X} \succeq m{0}$

 $\mathsf{rank}(\pmb{X}) = \mathsf{tr}(\pmb{X}) = k$

plus other combinatorial constraints

Ignoring rank constraint and relaxing combinatorial constraints on \boldsymbol{X} gives the semidefinite program:

$$\begin{array}{ll} \max & \operatorname{tr}(\boldsymbol{W}\boldsymbol{X}) \\ \text{s.t.} & \boldsymbol{X}\boldsymbol{e} \leq \boldsymbol{e} \\ & \operatorname{tr}(\boldsymbol{X}) = k \\ & \boldsymbol{X} \geq \boldsymbol{0}, \quad \boldsymbol{X} \succeq \boldsymbol{0}. \end{array}$$

Question: When does the optimal solution of this relaxation recover underlying cluster structure in similarity graph?

The Stochastic Block Model

Stochastic Block Model (SBM): generate random graph containing *k* clusters of size *r*:

- edges within clusters are added independently with probability \ensuremath{p}
- edges between-clusters are added with probability *q* < *p*.



Chen/Xu (2014) characterize when graphs sampled from the SBM are:

- trivial to cluster,
- easy to cluster (have polynomial-time algorithm),
- hard to cluster (via NP-hard max likelihood estimation)
- **impossible** to cluster (data has no meaningful cluster structure).

An *n*-node graph sampled from SBM is easy to cluster if

$$\frac{(p-q)^2}{q(1-q)} = \Omega\left(\frac{n}{r^2}\right).$$

Example: Clustered Euclidean data

Suppose each data point in the *i*th cluster C_i is placed uniformly at random in a ball centered at $c_i \in \mathbf{R}^d$.

Distance within clusters will be small compared to the distance between clusters if centers are well-separated.

Choose $w_{ij} = \exp(-\|\mathbf{x}^i - \mathbf{x}^j\|^2)$.





DOES NOT FIT STOCHASTIC BLOCK MODEL!!

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The Heterogeneous Planted Cluster Model

Assume that each node belongs to one of k clusters C_1, C_2, \ldots, C_k .

For each $u \in C_i$ and $v \in C_j$ we sample edge weight $w_{uv} = w_{vu}$ from distribution Ω_{ij} with

$$\boldsymbol{E}[w_{uv}] = \mu_{ij} \qquad \text{Var}[w_{uv}] = \sigma_{ij}^2 \qquad 0 \le w_{uv} \le 1.$$

Weights within the same block are i.i.d., but weight might not be identically distributed across blocks.



In the stochastic block model, we have perfect recovery if the gap constant $\gamma = q - p$ is sufficiently large.

In the heterogeneous case, we have perfect recovery if the **weak** assortativity constant

$$\gamma = \min_{\substack{q,s=1,2,\dots,k\\q\neq s}} \left\{ \mu_{qq} - \mu_{qs} \right\}$$

is sufficiently large.

Theorem (Pirinen-Ames 2019) Let $\hat{\sigma} := \max_{q \sigma_{qq}} \text{ and } \tilde{\sigma} := \max_{q,s} \sigma_{q,s}$.

Let \hat{r} denote size of the **smallest** planted cluster and r_{k+1} denote the number of outlier nodes.

Then there exists constant c > 0 such that if

$$\begin{split} \gamma \hat{r} &\geq c \max \Big\{ \sqrt{\tilde{\sigma}^2 n}, \sqrt{\tilde{\sigma}^2 \hat{r} \log n}, \sqrt{\hat{\sigma}^2 k r_{k+1}}, \\ &\sqrt{k r_{k+1} \log n / \hat{r}}, \ \mu_{k+1,k+1} r_{k+1}, \ \log n \Big\}. \end{split}$$

then we have perfect recovery with high probability.

Signal-to-noise ratio

Suppose that the edge weight is homogeneous: $\alpha = \mu_{qq}$, $\beta = \mu_{qs}$ for all $q \neq s$.

We can recover the planted clusters w.h.p. if

$$\frac{(\alpha-\beta)^2}{\tilde{\sigma}^2} = \Omega\left(\frac{n}{\hat{r}^2}\right).$$

The left-hand side acts as a signal-to-noise ratio: ratio of difference between expected edge weights to noise variance.

This agrees with/generalizes the easy regime for cluster recovery proposed by Chen and Xu (2014), and Jalali et al. (2015).

The relaxation is mostly parameter free: SDP needs number of clusters k but doesn't need estimate of cluster sizes r_i , gap statistic $\alpha - \beta$, etc., seen in similar theoretical guarantees.

Suppose Ω_1 and Ω_2 are Bernoulli distributions with probability of adding an edge p and q respectively (p > q) with no outliers $(r_{k+1} = 0)$.

Dense case: p, q constant (independent of n).

Have exact recovery w.h.p. if $\hat{r} \ge \hat{c}\sqrt{n}$ for some scalar \hat{c} (depending on p, q).

Sparse case: p constant, $q \le \frac{\log n}{n}$. Have exact recovery w.h.p. if $\hat{r} \ge \tilde{c} \log n$ for some constant \tilde{c} .

Rehnquist Supreme Court

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• Solve KDC with k = 2 to get the following partition of the Supreme court:

1: "Liberal"	2: "Conservative"
Stevens (St)	O'Connor (Oc)
Ginsberg (Gi)	Rehnquist (Re)
Souter (So)	Scalia (Sc) Thomas (Th)

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- Algorithm is sensitive to choice of *k*.
- Solve with k = 3:

Cluster 1	Cluster 2	Cluster 3
Thomas (Th) Scalia (Sc)	O'Connor (Oc) Kennedy (Ke) Rehnquist (Re)	Stevens (St) Breyer (Br) Ginsberg (Gi) Souter (So)

Current projects: Generalization of SBMs

More realistic planted models are needed:

- Overlapping clusters/communities;
- Finding largest of several planted clusters, possibly overlapping (without finding **all** clusters);
- Random graphs with **dependent** edges;
- Time-varying graphs; etc.,

Most machine learning **algorithms** are actually **heuristics**.

Approximately solve model problem for learning task (usually non-convex) and use approximate solution for inference process.

Would be extremely beneficial to have better understanding of the structure of local optima and optimization landscape of these model problems.

- Would allow better choices of initial solutions and heuristic parameters.
- Would encourage greater public trust in methods, more interpretability of results/predictions, etc.

Compressed sensing / LASSO: can find **sparsest** solution of underdetermined linear system by solving convex relaxation

 $\min\{\|\boldsymbol{x}\|_1: \boldsymbol{A}\boldsymbol{x} = \boldsymbol{b}\},\$

where $\|\mathbf{x}\|_1 = |x_1| + |x_2| + \cdots + |x_n|$, under certain assumptions about \mathbf{A} .

Rank minimization: can find **minimum rank** solution of linear system $\mathcal{A}(\mathbf{X}) = \mathbf{b}$ by solving

 $\min\{\|\boldsymbol{X}\|_*: \mathcal{A}(\boldsymbol{X}) = \boldsymbol{b}\},\$

under certain assumptions about \mathcal{A} , where $\|\boldsymbol{X}\|_*$ is the matrix nuclear norm.

Maximum Clique Problem: Ames/Vavasis 2011 showed that the maximum clique of graph G = (V, E) can be found by solving the relaxation

$$\min\left\{\|\boldsymbol{X}\|_*:\sum_{ij}x_{ij}=k,\ x_{ij}=0\ \forall\ ij\notin E\right\}$$

if *G* sampled from **planted clique model**. Recovery guarantee improved in Bombina/Ames 2020.

Similar average case recovery guarantees exist for **sparse PCA**, **nonnegative matrix factorization**, among other NP-hard problems.

Given set of objects and features, **biclustering** or **co-clustering** aims to partition both simultaneously so objects in bicluster strongly exhibit same features.

Want to obtain groups of objects similar with respect to a particular subset of features, while simultaneously grouping features.

Applications:

- identifying subsets of genes exhibiting similar expression patterns across subsets of experimental conditions in analysis of gene expression data,
- grouping documents by topics in document clustering, and
- grouping customers according to their preferences in collaborative filtering and recommender systems, etc.

The Biclustering SDP

Model the problem as **densest** *k*-**disjoint biclique problem**.

Let G = ((U, V), E) be a bipartite graph. Want collection of *k*-densest bipartite subgraphs, corresponding to *k* biclusters.



 $\begin{array}{ll} \max \ \operatorname{tr}(\boldsymbol{W}\boldsymbol{Z}) \\ \text{s.t.} \ \boldsymbol{Z}_{U,U}\boldsymbol{e} \leq \boldsymbol{e}, \quad \boldsymbol{Z}_{V,V}\boldsymbol{e} \leq \boldsymbol{e} \\ & \operatorname{tr}(\boldsymbol{Z}_{U,U}) = k = \operatorname{tr}(\boldsymbol{Z}_{V,V}) \\ & \boldsymbol{Z} \geq 0, \ \boldsymbol{Z} \in \boldsymbol{\Sigma}_{+}^{|\boldsymbol{U}|+|\boldsymbol{V}|} \end{array}$

Ames 2014 establishes conditions for perfect recovery in dense homogeneous case.

Would like to generalize to sparse heterogeneous case.

Current state of the art for solving clustering SDP requires $O(n^3)$ floating point operations for singular value decomposition each iteration; algorithm converges linearly.

Cannot solve large-scale problem instances.

Investigating intermediate relaxation via non-convex quadratic programming (QP).

Solve QP using linearized ADMM, with much lower iteration complexity.

- When do we have perfect recovery?
- When does our algorithm converge?

Current projects: Applied Data Analysis

Ford et al. 2021: applied novel classification algorithm to identify comprehension of language via EEG.

SLINGSHOT: machine learning pipeline for detecting lensed galaxies from observational telescope data.

Hyperspectral segmentation: Remote sensing (land-cover classification), biomedical samples (malignant vs. benign cells), geological samples (compositional/chronometric analysis)



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Software available from bpames.people.ua.edu/software