A constrained $\ell 1$ minimization approach for estimating multiple Sparse Gaussian or Nonparanormal Graphical Models

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Motivation: Structure Learning from Heterogeneous Samples

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- Learning relational graph structure among features/variables from an observed sample dataset is an important task in Machine Learning.
- This paper focuses on inferring graph structures from multiple related datasets (heterogeneous samples) about the same set of variables.
- We mainly focus on estimating conditional dependency graphs using the sparse Gaussian Graphical Model (sGGM).



Context/Task(1)

When Working on Multiple Different but Related Datasets:

- Samples of many real applications take the form of multiple different but related data matrices.
 - Blood cancer samples vs. Breast cancer samples;
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- Samples of many real applications take the form of multiple different but related data matrices.
 - Blood cancer samples vs. Breast cancer samples;
 - Normal patient samples vs. Cancel patient samples;
- A multi-task learning setting: to investigate the commonalities and differences among different datasets.



Our Aim: Shared and Task-specific Graph Structures

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Our Aim: Shared and Task-specific Graph Structures

- We aim to obtain shared and task-specific graph structures from heterogeneous samples.
- For example, in computational biology [Ideker and Krogan(2012)] urges to estimate housekeeping interactions and differential network among genes or proteins.



Our Aim: To Learn Shared and Task-specific Graph Structures from Multiple Related Datasets

Main Task: How to estimate / learn shared (Ω₅) and task-specific (Ω_l⁽ⁱ⁾) graph structures among feature variables from multiple different but related datasets about the same set of features.



Context/Task(2)



- X Data matrix.
- Σ Covariance matrix.
- Ω Inverse of covariance matrix (precision matrix).
- $X^{(i)}$ The *i*-th data matrix.
- $\Sigma^{(i)}$ The *i*-th covariance matrix.
- $\Omega^{(i)}$ The *i*-th precision matrix.
 - *p* The total number of feature variables.
 - n_i The number of samples in the *i*-th data matrix.
 - K The total number of tasks.

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- Precision matrix Ω is the inverse of covariance matrix Σ
- The sparsity pattern of Ω captures the conditional dependency pattern among variables.
- For example,



• Traditionally, we estimate sGGM from samples (of a single task) using an ℓ_1 penalized MLE formulation.

Graphical Lasso
[Friedman et al.(2008)Friedman, Hastie, and Tibshirani]

$$\underset{\Omega}{\operatorname{argmin}} - \ln \det(\Omega) + \operatorname{tr} \left(\Omega \widehat{\Sigma} \right) + \lambda_n ||\Omega||_1$$
(1.1)

Previous Methods: Joint Graphical Lasso (JGL) for Jointly Estimating Multiple sGGMs

• Most previous studies add a second penalty function *P*() into the penalized likelihood formulation.

Joint Graphical Lasso (JGL) [Danaher et al.(2013)Danaher, Wang, and Witten]

$$\underset{\Omega^{(i)}}{\operatorname{argmin}} - \sum_{i} n_{i} (\ln \det(\Omega^{(i)}) + \operatorname{tr}\left(\Omega^{(i)}\widehat{\Sigma}^{(i)}\right)) + \lambda_{1} \sum_{i} ||\Omega^{(i)}||_{1} + \lambda_{2} P(\Omega^{(1)}, \Omega^{(2)}, \dots, \Omega^{(K)})$$

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- P(Ω⁽¹⁾, Ω⁽²⁾,..., Ω^(K)) captures a certain assumption about relationships between multiple graphs.
- For example, fused norm to push graphs similar: $P(\Omega^{(1)}, \Omega^{(2)}, \dots, \Omega^{(K)}) = \sum_{i>j} ||\Omega^{(i)} - \Omega^{(j)}||_1.$

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- Two possible ways to infer multiple sGGMs from heterogeneous samples:
 - (1) Estimating one by one using graphical lasso by assuming the graphs are not related.
 - (2) Using JGL: joint graphical lasso by designing the appropriate second penalty function *P*().

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Drawbacks:

- I: Both of them can not directly output the shared structure among multiple graphs.
- II: Need extra steps to decode and can not control estimating the shared and task-specific pattern among graphs.
- **III:** No theoretical analysis in the previous JGL studies to prove why jointly learning graphs is helpful?

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- It can control the estimation of shared versus the task-specific patterns.
- It provides a strong theoretical guarantee.
- It achieves good empirical performance.

Proposed Method: Our "SIMULE" Formulation

We model each task's precision matrix $\Omega^{(i)}$ as a sum of task-specific $\Omega_I^{(i)}$ and task-shared Ω_S :

$$\Omega^{(i)} = \Omega_I^{(i)} + \Omega_S \tag{2.1}$$



Proposed method: Overview Figure



Why JGL Estimators Can't Get "SIMULE"

• JGL estimators are mostly solved by ADMM based optimization.

CLIME estimator [Cai et al.(2011)Cai, Liu, and Luo]

$$\begin{aligned} \underset{\Omega}{\operatorname{argmin}} & \|\Omega\|_{1} \\ \text{Subject to: } & \|\widehat{\Sigma}\Omega - I\|_{\infty} \leq \lambda_{n} \end{aligned} \tag{2.2}$$

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- With "SIMULE" formulation, difficult to separate the optimization into independent ADMM sub-procedures. Because,
 - The derivative of "SIMULE" in the JGL, i.e., gradient of $\ln \det(\Omega_I^{(i)} + \Omega_S)$ gets inverse of matrix summation.
 - Inverse of the summation of two matrices makes the optimization not separable.

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 - Inverse of the summation of two matrices makes the optimization not separable.
- Therefore, we use an alternative formulation for sGGM: A constrained ℓ_1 minimization formulation.

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SIMULE: to Infer <u>Shared and Individual Parts of MUL</u>tiple sGGM <u>Explicitly</u>

 By using a constrained l₁ minimization formulation, our estimator SIMULE can jointly learn multiple graphs from multiple different but related sample datasets (on the same set of feature variables).

SIMULE

$$\begin{split} \widehat{\Omega}_{I}^{(1)}, \widehat{\Omega}_{I}^{(2)}, \dots, \widehat{\Omega}_{I}^{(K)}, \widehat{\Omega}_{S} &= \underset{\Omega_{I}^{(i)}, \Omega_{S}}{\operatorname{argmin}} \sum_{i} ||\Omega_{I}^{(i)}||_{1} + \epsilon K ||\Omega_{S}||_{1} \end{split}$$

$$(2.3)$$
Subject to: $||\widehat{\Sigma}^{(i)}(\Omega_{I}^{(i)} + \Omega_{S}) - I||_{\infty} \leq \lambda_{n}, \ i = 1, \dots, K$

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- Can be solved by any linear programming solver.
- We have proved the "SIMULE" formulation guarantees a unique optimal solution.
- We use ϵ to control the sparsity of shared versus task-specific graph patterns.

Model Variation: NSIMULE for jointly estimating multiple nonparanormal Graphical Models

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- The only necessary change: by simply replacing the sample covariance matrices $\widehat{\Sigma}^{(i)}$ in Equation 2.3 into the kendal's tau correlation matrices $\widehat{\mathbf{S}}^{(i)}$.

Model Variation: NSIMULE for jointly estimating multiple nonparanormal Graphical Models

- The Gaussian assumption of our model can extend easily to a more general distribution family: nonparanormal.
- The only necessary change: by simply replacing the sample covariance matrices $\widehat{\Sigma}^{(i)}$ in Equation 2.3 into the kendal's tau correlation matrices $\widehat{\mathbf{S}}^{(i)}$.
- We denote this estimator as nonparanormal SIMULE (NSIMULE).

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• Comparing SIMULE v. CLIME w.r.t the statistical convergence rate for estimating *K* graphs:

Multi-task:K Single-task:
$$O(\frac{\log(Kp)}{n_{tot}})$$
 $\sum_{i} O(\frac{\log p}{n_i}))$

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- By assuming $n_i = \frac{n_{tot}}{K}$:
- We can conclude that $\frac{\log(Kp)}{n_{tot}} < K \frac{\log p}{n_{tot}}$
- This indicates that the multi-task estimator is better!!!

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Results on Synthetic Datasets: Accuracy and Parallelization

• Accuracy (AUC with a varying p and a varying K):



• Computation time cost with a varying *p*:



Results on Synthetic Datasets: Sensitivity of Hyperparameter ϵ

• The hyperpara ϵ controls the differences of sparsity among the shared graph and task-specific graphs.



• The sensitivity of ϵ vs. accuracy.



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Results on Two Real-World Datasets: Number of Matched Edges versus the Existing Domain Databases

- Two real world datasets:
 - (1) Gene expressions of samples in 2 different cell types
 - (2) Transcription Factors' ENCODE ChIP-seq measurements across 3 different cell lines
- Validation by counting the overlapped interactions according to the existing bio-databases (MInact).
- Our methods obtain the most matches compared to the state-of-the-art baselines.



- The project website: http://jointggm.org/
- R package "simule":
 - install.packages("simule")
 - demo(simuleDemo) !
 - https:

//cran.r-project.org/web/packages/simule/index.html

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