Learning Local Dependence In Ordered Data

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Abstract

In many applications, data come with a natural ordering. This ordering can often induce local dependence among nearby variables. However, in complex data, the width of this dependence may vary, making simple assumptions such as a constant neighborhood size unrealistic. We propose a framework for learning this local dependence based on estimating the inverse of the Cholesky factor of the covariance matrix. Penalized maximum likelihood estimation of this matrix yields a simple regression interpretation for local dependence in which variables are predicted by their neighbors. Our proposed method involves solving a convex, penalized Gaussian likelihood problem with a hierarchical group lasso penalty. The problem decomposes into independent subproblems which can be solved efficiently in parallel using first-order methods. Our method yields a sparse, symmetric, positive definite estimator of the precision matrix, encoding a Gaussian graphical model. We derive theoretical results not found in existing methods attaining this structure. In particular, our conditions for signed support recovery and estimation consistency rates in multiple norms are as mild as those in a regression problem. Empirical results show our method performing favorably compared to existing methods. We apply our method to genomic data to flexibly model linkage disequilibrium. Our method is also applied to improve the performance of discriminant analysis in sound recording classification.

1 Introduction

Estimating large inverse covariance matrices is a fundamental problem in modern multivariate statistics. Consider a random vector $X = (X_1, \ldots, X_p)^T \in \mathbb{R}^p$ with mean zero and covariance matrix $E(XX^T) = \Sigma$. Unlike the covariance matrix, which captures marginal correlations among variables in $X$, the inverse covariance matrix $\Omega = \Sigma^{-1}$ (also known as the precision matrix) characterizes conditional correlations and, under a Gaussian model, $\Omega_{jk} = 0$ implies that $X_j$ and $X_k$ are conditionally independent given all other variables. When $p$ is large, it is common to regularize the precision matrix estimator by making it sparse (see, e.g., Pourahmadi 2013). This paper focuses on the special context in which variables have a natural ordering, such as when data are collected over time or along a genome. In such a context, it is often reasonable to assume that random variables that are far away in the ordering are less dependent than those that are close together. For example, it is known that genetic mutations that occur close together on a chromosome are more likely to be co-inherited than mutations that are located far apart. We propose a method for estimating the precision matrix based on this assumption while also allowing each random variable to have its own notion of closeness.

In general settings where variables do not necessarily have a known ordering, two main types of convex methods with strong theoretical results have been developed to introduce sparsity in $\Omega$. The first approach, known as the graphical lasso (Yuan & Lin 2007, Banerjee et al. 2008, Friedman et al. 2008, Rothman et al. 2008), performs penalized maximum likelihood, solving
\[
\min_{\Omega \succ 0, \Omega = \Omega^T} \mathcal{L}(\Omega) + \lambda P(\Omega), \quad \text{where } \mathcal{L}(\Omega) = -\log \det \Omega + n^{-1} \sum_{i=1}^n x_i^T \Omega x_i \text{ is the negative log-likelihood of a sample of } n \text{ independent Gaussian random vectors and } P(\Omega) \text{ is the (vector) } \ell_1\text{-norm of } \Omega. \]

\[\text{Zhang & Zou (2014) generalizes this framework by introducing a new convex loss function called } D\text{-trace loss, and proposes a positive definite precision matrix estimator by minimizing a lasso-penalized version of this loss. The second approach is through penalized pseudo-likelihood, the most well-known of which is called neighborhood selection (Meinshausen & Bühlmann 2006). Estimators in this category are usually solved by a column-by-column approach and thus are more amenable to theoretical analysis (Yuan 2010, Cai et al. 2011, Liu & Luo 2012, Liu & Wang 2012, Sun & Zhang 2013, Khare & Zou 2014). However they are not guaranteed to be positive definite and do not exploit the symmetry of } \Omega. \]

\[\text{Peng et al. (2012) propose a partial correlation matrix estimator that develops a symmetric version of neighborhood selection; however, positive definiteness is still not guaranteed.} \]

In the context of variables with a natural ordering, by contrast, almost no work uses convex optimization to flexibly estimate \( \Omega \) while exploiting the ordering structure. Sparsity is usually induced via the Cholesky decomposition of \( \Sigma \), which leads to a natural interpretation of sparsity. Consider the Cholesky decomposition \( \Sigma = QQ^T \), which implies \( \Omega = L^T L \) for \( L = Q^{-1} \) for lower triangular matrices \( Q \) and \( L \) with positive diagonals. The assumption that \( X \sim N(0, \Sigma) \) is then equivalent to a set of linear models in terms of rows of \( L \), i.e., \( L_{11}X_1 = \varepsilon_1 \) and

\[
L_{rr}X_r = -\sum_{k=1}^{r-1} L_{rk}X_k + \varepsilon_r, \quad r = 2, \ldots, p, \tag{1}
\]

where \( \varepsilon \sim N(0, I_p) \). Thus, \( L_{rk} = 0 \) can be interpreted as saying that in predicting \( X_r \) from the previous random variables, one does not need to know \( X_k \). This observation has motivated previous work, including Pourahmadi (1999), Wu & Pourahmadi (2003), Huang et al. (2006), Shojaie & Michailidis (2010), Khare et al. (2016). While these methods assume sparsity in \( L \), they do not require local dependence, because each variable is allowed to be dependent on predecessors that are distant from it (compare the upper left to the upper right panel of Figure 10).

The assumption of “local dependence”, which unlike other methods, is our target of learning, can be expressed as saying that each variable \( X_r \) can be best explained by exactly its \( K_r \) closest predecessors:

\[
L_{rr}X_r = -\sum_{k=r-K_r}^{r-1} L_{rk}X_k + \varepsilon_r, \quad \text{for } L_{rk} \neq 0, \quad r - K_r \leq k \leq r - 1, \quad r = 2, \ldots, p. \tag{2}
\]

Note that this does not describe all patterns of a variable depending on its nearby variables. For example, \( X_r \) can be dependent on \( X_{r-2} \) but not on \( X_{r-1} \). In this case, the dependence is still local, but would not be captured by (2). Still we focus on the restricted class (2), since it greatly simplifies the interpretation of the learned dependence structure by capturing the extend of this dependence in a single number \( K_r \), the neighborhood size.

Bickel & Levina (2008) studied theoretical properties in the case that all bandwidths, \( K_r \), are equal, in which case model (2) is a \( K_r \)-ordered antedependence model (Zimmerman & Nunez-Anton 2009). A banded estimate of \( L \) then induces a banded estimate of \( \Omega \). The nested lasso approach of Levina et al. (2008) provides for “adaptive banding”, allowing \( K_r \) to vary with \( r \) (which corresponds to variable-order antedependence models in Zimmerman & Nunez-Anton 2009); however, their formulation is non-convex, meaning that their proposed algorithm does not necessarily minimize the stated objective and no theoretical properties of this estimator have been established.

In this paper, we propose a penalized likelihood approach that provides the flexibility of the nested lasso but is formulated as a convex optimization problem, which allows us to prove strong theoretical properties and to provide an efficient, scalable algorithm for computing the
estimator. The theoretical development of our method allows us to make clear comparisons with known results for the graphical lasso (Ravikumar et al. 2011) in the non-ordered case. Both methods are convex penalized likelihood approaches, so this comparison highlights the similarities and differences between the ordered and non-ordered problems.

There are two key choices we make that lead to a convex formulation. First, we express the optimization problem in terms of the Cholesky factor $L$. The nested lasso and other methods (starting with Pourahmadi (1999) use the modified Cholesky decomposition, $\Omega = T^T D^{-1} T$, where $T$ is a lower-triangular matrix with ones on its diagonal and $D$ is a diagonal matrix with positive entries. While $L(\Omega)$ is convex in $\Omega$, the negative log-likelihood $L(T^T D^{-1} T)$ is not jointly convex in $T$ and $D$. By contrast,

$$L(L^T L) = -\log \det (L^T L) + \frac{1}{n} \sum_{i=1}^{n} x_i^T L^T L x_i = -2 \sum_{r=1}^{p} \log L_{rr} + \frac{1}{n} \sum_{i=1}^{n} \|Lx_i\|^2_2$$  (3)

is convex in $L$. This parametrization is considered in Aragam & Zhou (2015), Khare et al. (2014) and Khare et al. (2016). Maximum likelihood estimation of $L$ preserves the regression interpretation by noting that

$$L(L^T L) = -2 \sum_{r=1}^{p} \log L_{rr} + \frac{1}{n} \sum_{r=1}^{p} \sum_{i=1}^{n} L_{rr}^2 \left( x_{ir} + \sum_{k<r} L_{rk} x_{ik} / L_{rr} \right)^2 .$$

This connection has motivated previous work with the modified Cholesky decomposition, in which $T_{rk} = -L_{rk} / L_{rr}$ are the coefficients of a linear model in which $X_r$ is regressed on its predecessors, and $D_{rr} = L_{rr}^{-2}$ corresponds to the residual variance. The second key choice is our use of a hierarchical group lasso in place of the nested lasso’s nonconvex penalty.

## 2 Estimator

For a given tuning parameter $\lambda \geq 0$, we define our estimator $\hat{L}$ to be a minimizer of the following penalized negative Gaussian log-likelihood

$$\hat{L} \in \arg \min_{L, L_{rr}\geq 0} \left\{ -2 \sum_{r=1}^{p} \log L_{rr} + \frac{1}{n} \sum_{i=1}^{n} \|Lx_i\|^2_2 + \lambda \sum_{r=2}^{p} P_r (L_r) \right\} ,$$  (4)

where the penalty for each row $P_r$ is defined by

$$P_r (L_r) = \sum_{\ell=1}^{r-1} \left\| W^{(\ell)} * L_{\ell,\ell} \right\|_2 = \sum_{\ell=1}^{r-1} \left( \sum_{m=1}^{\ell} w_{\ell m}^2 \right)^{1/2} .$$  (5)

The weight vector $W^{(\ell)} = (w_{\ell 1}, \ldots, w_{\ell \ell}) \in \mathbb{R}^\ell$, and $*$ is the element-wise multiplication. We denote $L_{\ell,\ell}$ to be a vector of elements of $L$ from the group $g_{\ell,\ell}$, which is defined as follows: for each row index $r$ and $1 \leq \ell \leq r-1$, the $\ell$-th group $g_{\ell,\ell} = \{(r, \ell') : \ell' \leq \ell\}$ corresponds to the first $\ell$ elements in the $r$-th row, and thus satisfies $g_{r,1} \subset g_{r,2} \subset \ldots \subset g_{r,r-1}$. Note that in (5), each row $r$ of $L$ is penalized with a sum of $r-1$ nested, weighted $L_2$-norm penalties. This is a hierarchical group lasso penalty (Yuan & Lin 2007, Zhao et al. 2009, Jenatton et al. 2011) with group structure conveyed in Figure 1.

With $w_{\ell m} > 0$, this nested structure always puts more penalty on those elements that are further away from the diagonal. Since the group lasso has the effect of setting to zero a subset of groups, it is apparent that this choice of groups ensures that whenever the elements in $g_{\ell,\ell}$ are set to zero, elements in $g_{\ell',\ell'}$ have already been set to zero for all $\ell' \leq \ell$. In other words, for each row of $\hat{L}$, the nonzeros are those elements within some (row-specific) distance of the diagonal.
This is in contrast to the $\ell_1$-penalty as used in [Khare et al. (2016)], which produces sparsity with no particular structure (compare top-left and top-right panel of Figure 10 for quantitative difference).

The choice of weights, $w_{\ell m}$, affects both the empirical and theoretical performance of the estimator. We focus primarily on a quadratically decaying set of weights,

$$w_{\ell m} = \frac{1}{(\ell - m + 1)^2}, \tag{6}$$

but also consider the unweighted case (in which $w_{\ell m} = 1$). The decay counteracts the fact that the elements of $L$ appear in differing numbers of groups (for example $L_{r1}$ appears in $r - 1$ groups whereas $L_{r,r-1}$ appears in just one group). In a related problem, [Bien et al. (2015)] choose weights that decay more slowly with $\ell - m$ than (6). Our choice makes the enforcement of hierarchy weaker so that our penalty behaves more closely to a non-hierarchical group lasso penalty, which, in our case, is a lasso penalty [Tibshirani 1996]. The choice of weight sequence in (6) is more amenable to theoretical analysis; however, in practice the unweighted case is more efficiently implemented and works well empirically.

![Figure 1: There are $\binom{d}{g}$ groups used in the penalty, with each row $r$ having $r - 1$ nested groups $g_{r,1} \subset g_{r,2} \subset \ldots \subset g_{r,r-1}$. Left: the group $g_{4,3}$. Middle: the nested group structure $g_{4,1} \subset g_{4,2} \subset g_{4,3}$. Right: A possible sparsity pattern in $\hat{L}$, where elements in $g_{2,1}, g_{4,2}$ (and thus $g_{4,1}$) and $g_{5,1}$ are set to zero.]

Problem (4) is convex in $L$. While $-\log \det(\cdot)$ is strictly convex, $- \sum g \log(L_{rr})$ is not strictly convex in $L$. Thus, the arg min in (4) may not be unique. In Section 4, we provide sufficient conditions to ensure uniqueness with high probability.

In Appendix A we show that (4) decouples into $p$ independent subproblems, each of which estimates one row of $L$. More specifically, let $X \in \mathbb{R}^{n \times p}$ be a sample matrix with rows $x_i \sim N(0, \Sigma)$, $\hat{L}_{11} = n^{1/2}(X_i^T X_i)^{-1/2}$ and for $r = 2, \ldots, p$,

$$\hat{L}_{r,1:r} = \arg\min_{\beta \in \mathbb{R}^p : \beta_r > 0} \left\{ -2 \log \beta_r + \frac{1}{n} \|x_{1:r} \|_2^2 + \lambda \sum_{\ell = 1}^{r-1} \left( \sum_{m=1}^{\ell} w_{\ell m}^2 \beta_m^2 \right)^{1/2} \right\}. \tag{7}$$

This observation means that the computation can be easily parallelized, which potentially can achieve a linear speed up with the number of CPU cores. Theoretically, to analyze the properties of $\hat{L}$ it is easier to start by studying an estimator of each row, i.e., a solution to (7). We will see in Section 4 that problem (7) has connections to a penalized regression problem, meaning that both the assumptions and results we can derive are better than working with simply estimating the matrix itself.

In light of the regression interpretation of (4), $\hat{L}$ provides an interpretable notion of local dependence; however, we can of course also use our estimate of $L$, to estimate $\Omega$: $\hat{\Omega} = \hat{L}^T \hat{L}$. By construction, this estimator is both symmetric and positive definite. Unlike a lasso penalty,
which would induce unstructured sparsity in the estimate of \( L \) and thus is not guaranteed to produce a sparse estimate of \( \Omega \), the adaptively banded structure in our estimator of \( L \) can yield a generally banded \( \hat{\Omega} \) (with bandwidth determined by the least sparse row of \( \hat{L} \). See the top-left and bottom-left panels in Figure 10 for an example).

3 Computation

As observed above, we can compute \( \hat{L} \) by solving (in parallel across \( r \)) problem \([7]\). Consider an alternating direction method of multiplier (ADMM) approach that solves the equivalent problem

\[
\min_{\beta, \gamma} \left\{ -2 \log \beta_r + \frac{1}{n} \|X_{1:r} \beta\|^2 + \lambda \sum_{\ell=1}^{r-1} \left( \sum_{m=1}^{\ell} w_{\ell m}^2 \gamma_{m}^2 \right)^{1/2} \right\} \quad \text{s.t. } \beta = \gamma
\]

Algorithm 1 presents the ADMM algorithm, which repeatedly minimizes this problem’s augmented Lagrangian over \( \beta \), then over \( \gamma \), and then updates the dual variable \( u \in \mathbb{R}^r \). The main computational effort in the algorithm is in solving \([8]\) and \([9]\). Note that \([8]\) has a smooth objective function. Straightforward calculus gives the closed-form solution (see Appendix B for detailed derivation),

\[
\beta_r^{(t+1)} = -\frac{B - \sqrt{B^2 - 8A}}{2A} > 0
\]

\[
\beta_{-r}^{(t+1)} = - \left( 2S_{-r,-r}^{(r)} + \rho I \right)^{-1} \left( 2S_{-r,r}^{(r)} \beta_r^{(t+1)} + u_{-r}^{(t)} - \rho \gamma_{-r}^{(t)} \right)
\]

where

\[
S^{(r)} = \frac{1}{n} X_{1:r}^T X_{1:r}
\]

\[
A = 4S_{-r,-r}^{(r)} \left( 2S_{-r,r}^{(r)} + \rho I \right)^{-1} S_{-r,r}^{(r)} - 2S_{r,r}^{(r)} - \rho < 0
\]

\[
B = 2S_{r,-r}^{(r)} \left( 2S_{-r,r}^{(r)} + \rho I \right)^{-1} \left( u_{-r}^{(t)} - \rho \gamma_{-r}^{(t)} \right) - u_r^{(t)} + \rho \gamma_r^{(t)}
\]
The closed-form update above involves matrix inversion. With \( \rho > 0 \), the matrix \( 2S^{(r)}_{-r,-r} + \rho I \) is invertible even when \( r > n \). Since determining a good choice for the ADMM parameter \( \rho \) is in general difficult, we adapt the dynamic \( \rho \) updating scheme described in Section 3.4.1 of [Boyd et al. 2011].

Solving (9) requires evaluating the proximal operator of the hierarchical group lasso with general weights. We adopt the strategy developed in [Bien et al. 2015] (based on a result of Jenatton et al. 2011), which solves the dual problem of (9) by performing Newton’s method on at most \( r - 1 \) univariate functions. The detailed implementation is given in Algorithm 3 in Appendix C. Each application of Newton’s method corresponds to performing an elliptical projection, which is a step of blockwise coordinate ascent on the dual of (9) (see Appendix D for details). Finally we observe in Algorithm 2 that for the unweighted case (\( w_{tm} = 1 \)), solving (9) is remarkably efficient.

**Algorithm 2** Algorithm for solving (9) for unweighted estimator

1. Initialize \( \gamma^{(t)} = \beta^{(t)} + u^{(t-1)} / \rho \) and \( \tau = \lambda / \rho \)
2. for \( \ell = 1, \ldots, r - 1 \) do
   \[
   \left( \gamma^{(t)} \right)_{1: \ell} = \left( 1 - \frac{\tau}{\left\| \left( \gamma^{(t)} \right)_{1: \ell} \right\|_2} \right) \left( \gamma^{(t)} \right)_{1: \ell}
   \]
3. return \( \gamma^{(t)} \).

4 Statistical properties

In this section we study the statistical properties of our estimator. We begin by introducing some notation used throughout the paper. For two sequences of constants \( a(n) \) and \( b(n) \), the notation \( a(n) = o(b(n)) \) means that for every \( \varepsilon > 0 \), there exists a constant \( N > 0 \) such that \( |a(n)/b(n)| \leq \varepsilon \) for all \( n \geq N \). And the notation \( a(n) = \mathcal{O}(b(n)) \) means that there exists a constant \( N > 0 \) and a constant \( M > 0 \) such that \( |a(n)/b(n)| \leq M \) for all \( n \geq N \). For a sequence of random variables \( A(n) \), the notation \( A(n) = \mathcal{O}_P(b(n)) \) means that for every \( \varepsilon > 0 \), there exists a constant \( M > 0 \) such that \( \mathbb{P}( |A(n)/b(n)| > M ) \leq \varepsilon \) for all \( n \).

For a vector \( v = (v_1, \ldots, v_p) \in \mathbb{R}^p \), we define \( \|v\|_1 = \sum_{j=1}^p |v_j| \), \( \|v\|_2 = (\sum_{j=1}^p v_j^2)^{1/2} \) and \( \|v\|_\infty = \max_j |v_j| \). For a matrix \( M \in \mathbb{R}^{n \times p} \), we define the element-wise norms by two vertical bars. Specifically, \( \|M\|_\infty = \max_{j,k} |M_{jk}| \) and Frobenius norm \( \|M\|_F = (\sum_{j,k} M_{jk}^2)^{1/2} \). For \( q \geq 1 \), we define the matrix-induced (operator) \( q \)-norm by three vertical bars: \( \|M\|_q = \max_{\|v\|_q = 1} \|Mv\|_q \). Important special cases include \( \|M\|_2 \), also known as the spectral norm, which is the largest singular value of \( M \), as well as \( \|M\|_1 = \max_k \sum_{j=1}^p |M_{jk}| \) and \( \|M\|_\infty = \max_j \sum_{k=1}^p |M_{jk}| \). Note that \( \|M\|_1 = \|M\|_\infty \) when \( M \) is symmetric. Given any two index sets \( T \) and \( T' \), denote the sub-vector \( v_T = (v_i)_{i \in T} \). We also let \( M_T \) be the \( p \times |T| \) matrix with columns selected from index set \( T \), and \( M_T^T \) be the \( |T| \times |T'| \) matrix with rows and columns of \( M \) indexed by \( T \) and \( T' \) respectively. Specifically, we use \( L_r \) to denote the \( r \)-th row of \( L \).

In what follows, we consider a lower triangular matrix \( L \) having row-specific bandwidths, \( K_r \). The first \( J_r = r - 1 - K_r \) elements of row \( r \) are zero and the band of nonzero off-diagonals (of size \( K_r \)) is denoted \( I_r = \{ J_r + 1, \ldots, r - 1 \} \). See Figure 2 for a graphical example of \( K_5, J_5, I_4 \) and \( T_4 \).
Figure 2: Schematic of the sparsity pattern assumption

Our theoretical analysis is built on the following assumptions:

**A1 Sparsity assumption:** The true Cholesky factor $L \in \mathbb{R}^{p \times p}$ is a lower triangular matrix such that the precision matrix $\Omega = L^T L$. $L$ has row-specific bandwidths $K_r$ such that $L_{rj} = 0$ for $0 < j < r - K_r$.

**A2 Gaussian assumption:** The sample matrix $X \in \mathbb{R}^{n \times p}$ has $n$ independent rows with each row $x_i$ drawn from $N(0, \Sigma)$.

**A3 Irrepresentable condition:** There exists some $\alpha \in (0, 1]$ such that
$$
\max_{2 \leq r \leq p} \max_{\ell \in I_r^c} \left\| \Sigma_{\ell \ell} (\Sigma_{\ell I_r})^{-1} \right\|_1 \leq \frac{6}{\pi^2} (1 - \alpha)
$$

**A4 Bounded singular values:** There exists a constant $\kappa$ such that
$$
0 < \kappa^{-1} \leq \sigma_{\text{min}}(L) \leq \sigma_{\text{max}}(L) \leq \kappa
$$

When $\max_r K_r < n$, the Gaussianity assumption [A2] implies that $X_{I_r}$ has full column rank for all $r$ with probability one. Our analysis applies to the general high-dimensional scaling scheme where $K_r = K_r(n)$ and $p = p(n)$ can grow with $n$.

For $r = 2, \ldots, p$ and $\ell \in I_r^c = \{1, \ldots, J_r, r\}$, let
$$
\theta^{(l)}_r := \text{Var} (X_\ell | X_{I_r}) \quad \text{and} \quad \theta_r := \max_{\ell \in I_r^c} \theta^{(l)}_r.
$$

By assumption [A2], $\theta^{(l)}_r = \Sigma_{\ell \ell} - \Sigma_{\ell I_r} (\Sigma_{I_r I_r})^{-1} \Sigma_{I_r \ell}$ represents the noise variance when regressing $X_\ell$ on $X_{I_r}$, i.e., for $\ell = 1, \ldots, J_r, r$,
$$
X_\ell = \Sigma_{\ell I_r} (\Sigma_{I_r I_r})^{-1} X_{I_r}^T + E_\ell \quad \text{with} \quad E_\ell \sim N \left(0, \theta^{(l)}_r\right).
$$

In words, $\theta^{(l)}_r$ measures the degree to which $X_\ell$ cannot be explained by the variables in the support and $\theta_r$ is the maximum such value over all $\ell$ outside of the support $I_r$ in the $r$-th row. Intuitively, the difficulty of the estimation problem increases with $\theta_r$. Note that for $r = 1, \ldots, p$, (11) implies $\theta^{(l)}_r = 1/L^2_{r r}$.

Assumption [A3] (along with the $\beta_{\text{min}}$ condition) is essentially a necessary and sufficient condition for the support recovery of the lasso-type methods (see, e.g., Zhao & Yu 2006, Meinshausen & Bühlmann 2006, Wainwright 2009, Van de Geer & Bühlmann 2009, Ravikumar et al.
The constant $\alpha \in (0, 1]$ is usually referred to as the irrepresentable (incoherence) constant. Intuitively, the irrepresentable condition requires low correlations between signals and noise predictors, and thus a value of $\alpha$ that is close to 1 implies that the support recovery is easier to achieve. The constant $6\pi^{-2}$ is determined by the choice of weight (6) and can be eliminated by absorbing its reciprocal into the definition of the weights $w_{\ell m}$. Doing so, one finds that our irrepresentable condition is essentially the same as the mild one found in the regression setting (Wainwright 2009) despite the fact that our goal is estimating a precision matrix.

Assumption A4 is a bounded singular value condition. By $\Omega = L^T L$, 

\begin{equation}
0 < \kappa^{-2} \leq \sigma_{\min}(\Sigma) \leq \sigma_{\max}(\Sigma) \leq \kappa^2 \tag{11}
\end{equation}

which is equivalent to the commonly used bounded eigenvalue condition in other literatures.

### 4.1 Row-specific results

We start by analyzing support recovery properties of our estimator for each row, i.e., the solution to the subproblem (7). For $r > n$, the Hessian of the negative log-likelihood is not positive definite, meaning that the objective function may not be strictly convex in $\beta$ and the solution not necessarily unique. Intuitively, if the tuning parameter $\lambda$ is large, the resulting row estimate $\hat{L}_r$ is sparse and thus includes most variation in a small subset of the $r$ variables. More specifically, for large $\lambda$, $\hat{I}_r \subseteq I_r$ and thus by assumption A2, $X_{\hat{I}_r}$ has full rank, which implies that $\hat{L}_r$ is unique. The series of technical lemmas in Appendix E precisely characterizes the solution.

The first part of the theorem below shows that with an appropriately chosen tuning parameter $\lambda$ the solution to (7) is sparse enough to be unique and that we will not over-estimate the true bandwidth. Knowing that the support of the unique row estimator $\hat{L}_r$ is contained in the true support reduces the dimension of the parameter space, and thus leads to a reasonable error bound. Of course, if our goal were simply to establish the uniqueness of $\hat{L}_r$ and that $\hat{K}_r = K_r$, we could trivially take $\lambda = \infty$ (resulting in $\hat{K}_r = 0$). The latter part of the theorem thus goes on to provide a choice of $\lambda$ that is sufficiently small to guarantee that $\hat{K}_r = K_r$ (and, furthermore, that the signs of all nonzeros are correctly recovered).

**Theorem 1.** Consider the family of tuning parameters

\begin{equation}
\lambda = \frac{8}{\alpha^2} \frac{\theta_r \log r}{n} \tag{12}
\end{equation}

and weights given by (6). Under assumptions A1–A4, if the tuple $(n, J_r, K_r)$ satisfies

\begin{equation}
n > \alpha^{-2} \left(3\pi^2 K_r + 8\right) \theta_r \kappa^2 \log J_r, \tag{13}
\end{equation}

then with probability greater than $1 - c_1 \exp \{-c_2 \min(K_r, \log J_r)\} - 7 \exp (-c_3 n)$ for some constants $c_1, c_2, c_3$ independent of $n$ and $J_r$, the following properties hold:

1. The row problem (7) has a unique solution $\hat{L}_r$ and $\hat{K}_r \leq K_r$.
2. The estimate $\hat{L}_r$ satisfies the element-wise $\ell_\infty$ bound,

\begin{equation}
\left\| \hat{L}_r - L_r \right\|_\infty \leq \lambda \left(4 \left\| (\Sigma_{I_r I_r})^{-1} \right\|_\infty + 5\kappa^2 \right). \tag{14}
\end{equation}

3. If in addition,

\begin{equation}
\min_{j \geq J_r + 1} |L_{rj}| > \lambda \left(4 \left\| (\Sigma_{I_r I_r})^{-1} \right\|_\infty + 5\kappa^2 \right), \tag{15}
\end{equation}

then exact signed support recovery holds: For all $j \leq r$, $\text{sign}(\hat{L}_{rj}) = \text{sign}(L_{rj})$. 

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In the classical setting where the ambient dimension $r$ is fixed and the sample size $n$ is allowed to go to infinity, $\lambda \to 0$ and the above scaling requirement is satisfied. By (14) the row estimator $\hat{L}_r$ is consistent as is the classical maximum likelihood estimator. Moreover, it recovers the true support since (15) holds automatically. In high-dimensional scaling, however, both $n$ and $r$ are allowed to change, and we are interested in the case where $r$ can grow much faster than $n$. Theorem 1 shows that, if $\| (\Sigma_{I_r I_r}^{-1})_{\infty} \| = O(1)$ and if $n$ can grow as fast as $K_r \log J_r$, then the row estimator $\hat{L}_r$ still recovers the exact support of $L_r$ when the signal is at least $O(\sqrt{\log r \over n})$ in size, and the estimation error $\max_j |\hat{L}_{rj} - L_{rj}|$ is $O(\sqrt{\log r \over n})$. Intuitively, for the row estimator to detect the true support, we require that the true signal be sufficiently large. The condition (15) imposes limitations on how fast the signal is allowed to decay, which is the analogue to the commonly known $\beta_{\min}$ condition in the support recovery of the lasso problem.

Remark 2. Both tuning parameter selection (12) and error bound (14) depend on the true covariance matrix via $\theta_r$. This quantity can be bounded by $\kappa^2$ as in (11). Using the fact that $(\Sigma_{I_r I_r}^{-1})_{\ell,\ell}$ is positive definite:

$$\theta_r = \max_{\ell \in I_r} \theta_r^{(\ell)} = \max_{\ell \in I_r} \left\{ \Sigma_{\ell\ell} - \Sigma_{\ell r} (\Sigma_{I_r I_r}^{-1})_{\ell r} \right\} \leq \max_{\ell \in I_r} \Sigma_{\ell\ell} \leq \kappa^2.$$ 

The proof of Theorem 1 shows that the results in this theorem still hold true if we replace $\theta_r$ by $\kappa^2$. This observation leads to the fact that we can select a tuning parameter having the properties of the theorem that does not depend on the unknown sparsity level $K_r$. Therefore, our estimator is adaptive to the underlying unknown bandwidths.

4.1.1 Connections to the regression setting

Recall that in (1) we showed that estimation of the $r$-th row of $L$ can be interpreted as a regression of $X_r$ on its predecessors. It is thus very interesting to compare Theorem 1 to the standard high-dimensional regression results. Consider the following linear model of a vector $\mathbf{y} \in \mathbb{R}^n$ of the form

$$\mathbf{y} = \mathbf{Z} \eta + \omega \quad \omega \sim N(0, \sigma^2 I_n)$$

where $\eta \in \mathbb{R}^p$ is the unknown but fixed parameter to estimate, $\mathbf{Z} \in \mathbb{R}^{n \times p}$ is the design matrix with each row an observation of $p$ predictors, $\sigma^2$ is the variance of the zero-mean additive noise $\omega$. A standard approach in the high-dimensional setting where $p \gg n$ is the lasso (Tibshirani 1996), which solves the following convex optimization problem,

$$\min_{\eta \in \mathbb{R}^p} \frac{1}{2n} \| \mathbf{y} - \mathbf{Z} \eta \|^2_2 + \lambda \| \eta \|_1,$$

where $\lambda > 0$ is a regularization parameter. In the setting where $\eta$ is assumed to be sparse, the lasso solution is known to be able to successfully recover the signed support of the true $\eta$ with high probability when $\lambda$ is of the scale $\sigma \sqrt{\log p \over n}$ and certain technical conditions are satisfied (Wainwright 2009).

Despite the added complications of working with the log term in the objective of (7), Theorem 1 gives a clear indication that, in terms of difficulty of support recovery, the row estimate problem (7) is essentially the same as a lasso problem with random design, i.e., with each row $z_i \sim N(0, \Sigma)$ (Theorem 3, Wainwright 2009). Indeed, a comparison shows that the two irrepresentable conditions are equivalent. Moreover, $\theta_r$ plays the same role as Wainwright (2009)’s $\max_i \left( \Sigma_{S^c \hbar^c} - \Sigma_{S S} (\Sigma_{S S})^{-1} \Sigma_{S S^c} \right)_{ii}$, a threshold constant of the conditional covariance, where $S$ is the support of the true $\eta$.  

Städtler et al. (2010) introduced an alternative approach to the lasso in the context of the penalized mixture regression models, which solves the following optimization problem,

\[
(\hat{\phi}, \hat{\rho}) = \arg \min_{\phi, \rho} -2 \log \rho + \frac{1}{n} \| \rho y + Z \phi \|_2^2 + \lambda \| \phi \|_1,
\]

(18)

and \( \hat{\sigma} = \hat{\rho}^{-1}, \hat{\eta} = \hat{\phi} / \hat{\rho} \). Note that (18) basically coincides with (7) except for the penalty.

In Städtler et al. (2010), authors studied the asymptotic and non-asymptotic properties of the \( \ell_1 \)-penalized estimator for the general mixture regression models where the loss functions are non-convex. The theoretical properties of (18) are studied in Sun & Zhang (2010), which partly motivates the scaled lasso (Sun & Zhang 2012).

The theoretical work of Sun & Zhang (2010) differs from ours both in that they study the \( \ell_1 \) penalty (instead of hierarchical group lasso) and in their assumptions. The nature of our problem requires the sample matrix to be random (as in A2), while Sun & Zhang (2010) considers the fixed design setting, which does not apply in our context. Moreover, they provide prediction consistency and a deviation bound of the regression parameters estimation in \( \ell_1 \) norm. We give exact signed support recovery results for the regression parameters as well as estimation deviation bounds in various norm criteria. Also, they take an asymptotic point of view while we give finite sample results.

4.2 Matrix bandwidth recovery result

With the properties of the row estimators in place, we are ready to state results about estimation of the matrix \( L \). The following theorem gives an analogue to Theorem 1 in the matrix setting. Under similar conditions, with one particular choice of tuning parameter, the estimator recovers the true bandwidth for all rows adaptively with high probability.

**Theorem 3.** Let \( \theta = \max_r \theta_r \) and \( K = \max_r K_r \), and take

\[
\lambda = \frac{8}{\alpha} \sqrt{\frac{2\theta \log p}{n}}
\]

(19)

and weights given by (6). Under assumptions A1–A4, if \((n, p, K)\) satisfies

\[
n > \alpha^{-2} \theta \kappa^2 (12\pi^2 K + 32) \log p,
\]

(20)

then with probability greater than \( 1 - cp^{-1} \) for some constant \( c \) independent of \( n \) and \( p \), the following properties hold:

1. The estimator \( \hat{L} \) is unique, and it is at least as sparse as \( L \), i.e., \( \hat{K}_r \leq K_r \) for all \( r \).

2. The estimator \( \hat{L} \) satisfies the element-wise \( \ell_\infty \) bound,

\[
\| \hat{L} - L \|_\infty \leq \lambda \left( 4 \max_r \left\| \left( \Sigma_{L_r, L_r} \right)^{-1} \right\|_\infty + 5\kappa^2 \right).
\]

(21)

3. If in addition,

\[
\min_r \min_{j \geq \ell_r + 1} |L_{rj}| > \lambda \left( 4 \max_r \left\| \left( \Sigma_{L_r, L_r} \right)^{-1} \right\|_\infty + 5\kappa^2 \right),
\]

(22)

then exact signed support recovery holds: \( \text{sign}(\hat{L}_{rj}) = \text{sign}(L_{rj}) \) for all \( r \) and \( j \).

As discussed in Remark 2, we can replace \( \theta \) with its upper bound \( \kappa^2 \), and the results remain true. This theorem shows that one can properly estimate the sparsity pattern across all rows exactly using only one tuning parameter chosen without any prior knowledge of the true
bandwidths. In Section 4.1.1, we noted that the conditions required for support recovery and element-wise $\ell_\infty$ error bound for estimating a row of $L$ is similar to those of the lasso in the regression setting. A union bound argument allows us to translate this into exact bandwidth recovery in the matrix setting and to derive a reasonable convergence rate under conditions as mild as that of a lasso problem with random design. This technique is similar in spirit to neighborhood selection (Meinshausen & Bühlmann 2006), though our approach is likelihood-based.

Comparing (20) to (13), we see that the sample size requirement for recovering $L$ is determined by the least sparse row. While intuitively one would expect the matrix problem to be harder than any single row problem, we see that in fact the two problems are basically of the same difficulty (up to a multiplicative constant).

In the setting where variables exhibit a natural ordering, Shojaie & Michailidis (2010) proposed a penalized likelihood framework like ours to estimate the structure of directed acyclic graphs (DAGs). Their method focuses on variables which are standardized to have unit variance. In this special case, penalized likelihood does not involve the log-determinant term and under similar assumptions to ours, they proved support recovery consistency. However, they use lasso and adaptive lasso (Zou 2006) penalties, which do not have the built-in notion of local dependence. Since these $\ell_1$-type penalties do not induce structured sparsity in the Cholesky factor, the resulting precision matrix estimate is not necessarily sparse. By contrast, our method does not assume unit variances and learns an adaptively banded structure for $\hat{L}$ that leads to a sparse $\hat{\Omega}$ (thereby encoding conditional dependencies).

To study the difference between the ordered and non-ordered problems, we compare our method with Ravikumar et al. (2011), who studied the graphical lasso estimator in a general setting where variables are not necessarily ordered. Let $S$ index the edges of the graph specified by $\Omega = \Sigma^{-1}$. The sparsity recovery result and convergence rate are established under an irrepresentable condition imposed on $\Gamma = \Sigma \otimes \Sigma \in \mathbb{R}^{p^2 \times p^2}$:

$$\max_{e \in S} \left\| \Gamma_{es} \left( \Gamma_{SS} \right)^{-1} \right\|_1 \leq (1 - \alpha)$$

(23)

for some $\alpha \in (0, 1]$. Our assumption $\text{A3}$ is on each variable through the entries of the true covariance $\Sigma$ while (23) imposes such a condition on the edge variables $Y_{j,k} = X_j X_k - E(X_j X_k)$, resulting in a vector $\ell_1$-norm restriction on a much larger matrix $\Gamma$, which can be more restrictive for large $p$. More specifically, condition (23) arises in Ravikumar et al. (2011) to tackle the analysis of the log det $\Omega$ term in the graphical lasso problem. By contrast, in our setting the parameterization in terms of $L$ means that the log det term is simply a sum of log terms on diagonal elements and is thus easier to deal with, leading to the milder irrepresentable assumption. Another difference is that they require the sample size $n > c \kappa^2 d^2 \log p$ for some constant $c$. The quantity $d$ measures the maximum number of nonzero elements in each row of the true $\Sigma$, which in our case is $2K + 1$, and $\kappa = \left\| \left( \Gamma_{SS} \right)^{-1} \right\|_{\infty}$ can be much larger than $\kappa^2$.

Thus, comparing to (20), one finds that their sample size requirement is much more restrictive. A similar comparison could also be made with the lasso penalized D-trace estimator (Zhang & Zou 2014), whose irrepresentability condition involves $\Gamma = (\Sigma \otimes I + I \otimes \Sigma) / 2 \in \mathbb{R}^{p^2 \times p^2}$. Of course, the results in both Ravikumar et al. (2011) and Zhang & Zou (2014) are invariant to the permutation among variables; additionally, the random vector only needs to satisfy an exponential-type tail condition.

### 4.3 Precision matrix estimation consistency

Although our primary target of interest is $L$, the parameterization $\Omega = L^T L$ makes it natural for us to try to connect our results of estimating $L$ with the vast literature in directly estimating $\Omega$, which is the standard estimation target when the known ordering is not available. In this
section, we consider the estimation consistency of $\Omega$ using the results we obtained for $L$. The following theorem gives results of how well $\hat{\Omega} = \hat{L}^T L$ performs in estimating the true precision matrix $\Omega = L^T L$ in terms of various matrix norm criteria.

**Theorem 4.** Let $\theta = \max_r \theta_r$, $K = \max_r K_r$ and $s = \sum_r K_r$ denote the total number of nonzero off-diagonal elements in $L$. Define $\zeta_\Sigma = 8\sqrt{2\alpha} \left( 4 \max_r \left( \left\| \left( \Sigma_r I_r \right)^{-1} \right\|_\infty + 5\kappa^2 \right) \right)$. Under the assumptions in Theorem 3 the following deviation bounds hold with probability greater than $1 - cp^{-1}$ for some constant $c$ independent of $n$ and $p$:

$$\left\| \hat{\Omega} - \Omega \right\|_\infty \leq 2\zeta_\Sigma \|L\|_\infty \sqrt{\frac{\log p}{n}} + \zeta_\Sigma^2 (K + 1) \frac{\log p}{n},$$

$$\left\| \hat{\Omega} - \Omega \right\|_\infty \leq 2\zeta_\Sigma \|L\|_\infty (K + 1) \sqrt{\frac{\log p}{n}} + \zeta_\Sigma^2 (K + 1)^2 \frac{\log p}{n},$$

$$\left\| \hat{\Omega} - \Omega \right\|_2 \leq 2\zeta_\Sigma \|L\|_\infty (K + 1) \sqrt{\frac{\log p}{n}} + \zeta_\Sigma^2 (K + 1)^2 \frac{\log p}{n},$$

$$\left\| \hat{\Omega} - \Omega \right\|_F \leq 2\zeta_\Sigma \sqrt{(s + p) \log p} + \zeta_\Sigma^2 (K + 1) \sqrt{s + p} \frac{\log p}{n}.$$

When the quantities $\zeta_\Gamma$, $\|L\|_\infty$, and $\kappa$ are treated as constants, these bounds can be summarized more succinctly as follows:

**Corollary 5.** Using the notation, and conditions in Theorem 4 if $\zeta_\Gamma$, $\|L\|_\infty$ and $\kappa$ remain constant, then the scaling $(K + 1)^2 \log p = o(n)$ is sufficient to guarantee the following estimation error bounds:

$$\left\| \hat{\Omega} - \Omega \right\|_\infty = O_P \left( \sqrt{\frac{\log p}{n}} \right),$$

$$\left\| \hat{\Omega} - \Omega \right\|_\infty = O_P \left( (K + 1) \sqrt{\frac{\log p}{n}} \right),$$

$$\left\| \hat{\Omega} - \Omega \right\|_2 = O_P \left( (K + 1) \frac{\log p}{n} \right),$$

$$\left\| \hat{\Omega} - \Omega \right\|_F = O_P \left( \sqrt{(s + p) \log p} \frac{\log p}{n} \right).$$

The conditions for these deviation bounds to hold are those required for support recovery as in Theorem 3. In many cases where estimation consistency is more of interest than support recovery, we can still deliver the desired error rate in Frobenius norm, matching the rate derived in Rothman et al. (2008). In particular, we can drop the strong irrepresentable assumption and weaken the Gaussian assumption to the following marginal sub-Gaussian assumption:

**A5 Marginal sub-Gaussian assumption:** The sample matrix $X \in \mathbb{R}^{n \times p}$ has $n$ independent rows with each row drawn from the distribution of a zero-mean random vector $X = (X_1, \cdots, X_p)^T$ with covariance $\Sigma$ and sub-Gaussian marginals, i.e.,

$$\mathbb{E}\exp\left( tX_j / \sqrt{\Sigma_{jj}} \right) \leq \exp \left( Ct^2 \right)$$

for all $j = 1, \cdots, p$, $t \geq 0$ and for some constant $C > 0$ that is independent of $j$.

**Theorem 6.** Under assumption A1, A4, and A5, with tuning parameter $\lambda$ of scale $\sqrt{\frac{\log p}{n}}$ and weights as in (6), the scaling $(s + p) \log p = o(n)$ is sufficient for the following estimation error
bounds in Frobenius norm to hold:

\[ \| \hat{L} - L \|_F = \mathcal{O}_P \left( \frac{(s + p) \log p}{n} \right), \]
\[ \| \hat{\Omega} - \Omega \|_F = \mathcal{O}_P \left( \frac{(s + p) \log p}{n} \right). \]

The rates in Corollary 3 (and Theorem 4) essentially match the rates obtained in methods that directly estimate \( \Omega \) (e.g., the graphical lasso estimator, studied in Rothman et al. 2008 and Ravikumar et al. 2011) and the column-by-column methods as in Cai et al. 2011, Liu & Wang 2012 and Sun & Zhang 2013. However, the exact comparison in rates with these methods is not straightforward. First, the targets of interest are different. In the setting where the variables have a known ordering, we are more interested in the structural information among variables that is expressed in \( L \), and thus accurate estimation of \( L \) is more important. When such ordering is not available as considered in Rothman et al. (2008), Cai et al. (2011), Liu & Wang (2012) and so on, however, the conditional dependence structure encoded by the sparsity pattern in \( \Omega \) is more of interest, and the accuracy of directly estimating \( \Omega \) is the focus. Moreover, deviation bounds of different methods are built upon assumptions that treat different quantities as constants. Quantities that are assumed to remain constant in the analysis of one method might actually be allowed to scale with ambient dimension in a nontrivial manner in another method, which makes direct rate comparison among different methods complicated and less illuminating.

Our analysis can be extended to the unweighted version of our estimator, i.e., with weight \( w_{\ell m} = 1 \), but under more restrictive conditions and with slower rates of convergence. Specifically, assumption \( \text{A3} \) becomes \( \max_{\ell \in I_r} \left\| \Sigma_{\ell \ell_r} (\Sigma_{I_r I_r})^{-1} \right\|_1 \leq (1 - \alpha) / K_r \) for each \( r = 2, \ldots, p \). With the same tuning parameter choice \( (12) \) and \( (19) \), the terms of \( K_r \) and \( K_r \) in sample size requirements \( (13) \) and \( (20) \) are replaced with \( K_r^2 \) and \( K_r^2 \) respectively. The estimation error bounds in all norms are multiplied by an extra factor of \( K_r \). All of the above indicates that in highly sparse situations (in which \( K \) is very small), the unweighted estimator has very similar theoretical performance to the weighted estimator.

5 Simulation study

In this section we study the empirical performance of our estimators (with weights as in (6) and with no weights, i.e., \( w_{\ell m} = 1 \)) in simulated data. For comparison, we include two other sparse precision matrix estimators designed for the ordered-variable case:

- **Non-Adaptive Banding (Bickel & Levina 2008):** The method estimates \( L \) as a lower-triangular matrix with a fixed bandwidth \( K \) applying across all rows. The regularization parameter used in this method is the fixed bandwidth \( K \).

- **Nested Lasso (Levina et al. 2008):** This method yields an adaptive banded structure by solving a set of penalized least square problems (both the loss function and the nested-lasso penalty are non-convex). The regularization parameter controls the amount of penalty and thus the sparsity level of the resulting estimate.

All simulations are run at a sample size of \( n = 100 \), where each sample is drawn independently from the \( p \)-dimensional normal distribution \( N(0, (L^T L)^{-1}) \). We compare the performance of our estimators with the methods above both in terms of support recovery (in Section 5.1) and in terms of how well \( \hat{L} \) estimates \( L \) (in Section 5.2). For support recovery, we consider \( p = 200 \) and for estimation accuracy, we consider \( p = 50, 100, 200 \), which corresponds to settings where \( p < n, p = n, \) and \( p > n \), respectively.
We simulate with the following models of $L$. For each model, we adapt the parameterization $L = D^{-1}T$ as in [Khare et al., 2016], where $D$ is a diagonal matrix with diagonal elements drawn randomly from a uniform distribution on the interval $[2,5]$, and $T$ is a lower-triangular matrix with ones on its diagonal and off-diagonal elements defined as follows:

- **Model 1**: At one extreme of bandedness of the Cholesky factor $L$, we take the lower triangular matrix $L \in \mathbb{R}^{p \times p}$ to have a strictly banded structure, with each row having the same bandwidth $K_r = K = 1$ for all $r$. Specifically, we take $T_{rr} = 1$, $T_{r,r-1} = 0.8$ and $T_{r,j} = 0$ for $j < r - 1$.

- **Model 2**: At the other extreme, we allow $K_r$ to vary with $r$. We take $T$ to be a block diagonal matrix with 5 blocks, each of size $p/5$. Within each block, with probability 0.5 each row $r$ is assigned with a non-zero bandwidth that is randomly drawn from a uniform distribution on $\{1, \ldots, r-1\}$ (for $r > 1$). Each non-zero element in $T$ is then drawn independently from a uniform distribution on the interval $[0.1, 0.4]$, and is assigned with a positive/negative sign with probability 0.5.

- **Model 3**: Model 3 is a denser and thus more challenging version of Model 2, with $T$ a block diagonal matrix with only 2 blocks. Each of the blocks is of size $p/2$ and is generated the same way as in Model 2.

- **Model 4**: Model 4 is a dense block diagonal model. The matrix $T$ has a totally dense lower-triangular block from the $p/4$-th row to the $3p/4$-th row and takes value of zero everywhere else. Within this block, all off-diagonal elements are drawn uniformly from $[0.1, 0.2]$, and positive/negative signs are then assigned with probability 0.5.

Model 1 is a stationary autoregressive model with order 1. By the regression interpretation [1], for each $r$, it can be verified that the autoregressive polynomial of the $r$-th row of Models 2, 3, and 4 has all roots outside the unit circle, which characterizes stationary autoregressive models with lags equal to the corresponding row-wise bandwidths. See Figure 3 for examples of the four sparsity patterns for $p = 100$. The non-adaptive banding method should benefit from Model 1 while the nested lasso and our estimators are expected to perform better in the other three models where each row has its own bandwidth.

![Figure 3: Schematic of four simulation scenarios with $p = 100$: (from left to right) Model 1 is strictly banded, Model 2 has small variable bandwidth, Model 3 has large variable bandwidth, and Model 4 is block-diagonal. Black, gray, and white stand for positive, negative, and zero entries respectively. The proportion of number of nonzeros are 4%, 6%, 15%, and 26% respectively.](image)

Note that for all four models and every value of $p$ considered, we verified that the assumptions $A3$ and $A4$ hold and then simulated $n = 100$ observations according to each of the four models based on assumption $A2$.
5.1 Support recovery

We first study how well the different estimators correctly identify zeros in the four models above. We generate $n = 100$ random samples from each model with $p = 200$. The tuning parameter $\lambda \geq 0$ in (4) measures the amount of regularization and determines the sparsity level of the estimator. We use 100 tuning parameter values for each estimator and repeat the simulation 10 times.

Figure 4 shows the sensitivity (fraction of true non-zeros that are correctly recovered) and specificity (fraction of true zeros that are correctly set to zero) of each method parameterized by its tuning parameter (in the case of the non-adaptive banding, the parameter is the bandwidth itself, ranging from 0 to $p-1$). Each set of 10 curves of the same color corresponds to the results of one estimator, and each curve within the set corresponds to the result of one draw from 10 simulations. Curves closer to the upper-right corner indicate better classification performance (the $x + y = 1$ line corresponds to random guessing).

The sparsity level of the non-adaptive banding estimator depends only on the pre-specified bandwidth (which is the method’s tuning parameter) and not on the data itself. Consequently, the sensitivity-specificity curves for the non-adaptive banding do not vary across replications when simulating from a particular underlying model. The sparsity levels of the nested lasso and our methods, by contrast, hinge on the data, thus giving a different curve for each replication.

In practice, we find that our methods and the nested lasso sometimes produce entries with
very small, but nonzero, absolute values. To study support recovery, we set all estimates whose absolute values below $10^{-10}$ to zero, both in our estimators and the nested lasso.

In Model 1, we observe that all methods considered attain perfect classification accuracy for some value of their tuning parameter. While the non-adaptive approach is guaranteed to do so in this scenario, it is reassuring to see that the more flexible methods can still perfectly recover this sparsity pattern.

In Model 2, we observe that our two methods outperform the nested lasso, which itself, as expected, outperforms the non-adaptive banding method. As the model becomes more challenging (from Model 2 to Model 4), the performances of all four methods start deteriorating. Interestingly, the nested lasso no longer retains its advantage over the non-adaptive banding in model 3 and 4, while the performance advantage of our methods become even more substantial.

The fact that the unweighted version of our method outperforms the weighted version stems from the fact that all models are comparatively sparse for $p = 200$, and so the heavier penalty on each row delivered by the unweighted approach recovers the support more easily than the weighted version.

### 5.2 Estimation accuracy

We proceed by comparing the estimators in terms of how far $\hat{L}$ is from $L$. To this end, we generate $n = 100$ random samples from the four models with $p = 50$, $p = 100$, and $p = 200$. Each method is computed with the best tuning parameter selected to maximize the Gaussian likelihood on the validation data in a 5-fold cross-validation. For comparison, we report the estimation accuracy of each estimate in the scaled Frobenius norm $\frac{1}{p} \|L - L\|_F^2$, matrix infinity norm $\|L - L\|_\infty$, spectral norm $\|L - L\|_2$ and the scaled Kullback-Leibler loss (see, e.g., [Levina et al. 2008]).

The simulation is repeated 50 times, and the results are summarized in Figure 5 through Figure 8. Each figure, representing the results of all four methods applied to data generated from one model, has a 3-by-4 panel layout, where each row displays results in box-plots for three different values of $p$, and each row contains the results (in box-plots) of the estimation accuracy evaluated in one of the four criteria.
Figure 5: Estimation accuracy when data are generated from Model 1 which is strictly banded.
Figure 6: Estimation accuracy when data are generated from Model 2 which has small variable bandwidth.
Figure 7: Estimation accuracy when data are generated from Model 3 which has large variable bandwidth.
As expected, the non-adaptive banding estimator does better than the other estimators in Model 1. In Models 2, 3, and 4, where bandwidths vary with row, our estimators and the nested lasso outperform the non-adaptive banding.

A similar pattern is observed as in support recovery. As the model becomes more complex and $p$ gets larger, the performance of the nested lasso degrades and gradually becomes worse than the non-adaptive banding. By contrast, as the estimation problem becomes more difficult, the advantage in performance of our methods becomes more obvious.

We again observe that the unweighted estimator performs better than the weighted one. As shown in Section 4, the overall performance of our method hinges on the underlying model complexity (measured in terms of $\max_r K_r$) as well as the relative size of $n$ and $p$. When $n$ is
relatively small, usually a more constrained method (like the unweighted estimator) is preferred over a more flexible method (like the weighted estimator). So in our simulation setting, it is reasonable to observe that the unweighted method works better. Note that as the underlying $L$ becomes denser (from Model 1 to Model 4), the performance difference between the weighted and the unweighted estimator diminishes. This verifies our discussion in the end of Section 4 that the theoretical performance of the unweighted estimator becomes worse when the underlying model is dense.

6 Applications to data examples

In this section, we illustrate the practical merits of our proposed method by applying it to two data examples. We start with an application to genomic data where our method can help model the local correlation along the genome. In Section 6.2 we compare our method with other estimators within the context of a sound recording classification problem.

6.1 An application to genomic data

We consider an application of our estimator to modeling correlation along the genome. Genetic mutations that occur close together on a chromosome are more likely to be co-inherited than mutations that are located far apart (or on separate chromosomes). This leads to local correlations between genetic variants in a population. Biologists refer to this local dependence as linkage disequilibrium (LD). The width of this dependence is known to vary along the genome due to the variable locations of recombination hotspots, which suggests that adaptively banded estimators may be quite suitable in these contexts.

We study HapMap phase 3 data from the International HapMap project (Consortium et al. 2010). The data consist of $n = 167$ humans from the YRI (Yoruba in Ibadan, Nigeria) population, and we focus on $p = 201$ consecutive tag SNPs on chromosome 22 (after filtering out infrequent sites with minor allele frequency $\leq 10\%$).

Note that tag SNP data, which takes discrete values $\{0, 1, 2\}$, are non-Gaussian. However we argue that our estimator is still sensible to use in this case. First, the parameterization $\Omega = L^T L$ does not depend on the Gaussian assumption. Moreover the estimator corresponds to minimizing a penalized Bregman divergence of the log-determinant function (Ravikumar et al. 2011). Actually, by looking at the minimization problem (4), we find that the least square term can be interpreted as minimizing the prediction error in the linear models (1) while the log terms act as log-barrier functions to impose positive diagonal entries (which ensures that the resulting $\hat{L}$ is a valid Cholesky factor).

To gauge the performance of our estimator on modeling LD, we randomly split the 167 samples into training and testing sets of sizes 84 and 83 respectively. Along a path of tuning parameters with decreasing values, estimators $\hat{L}$ are computed on the training data. To evaluate $\hat{L}$ on a vector $\tilde{x}$ from the test data set, we can compute the error in predicting $\hat{L}_{rr}\tilde{x}_r$ using $-\sum_{k=1}^{p-1} \hat{L}_{r,k} \tilde{x}_k$ via (1) for each $r$, giving the error

$$
\text{err}(\tilde{x}) = \frac{1}{p-1} \sum_{r=2}^{p} \left( \hat{L}_{rr}\tilde{x}_r + \sum_{k=1}^{r-1} \hat{L}_{rj}\tilde{x}_k \right)^2.
$$

This quantity (with mean and the standard deviation over test samples) is reported in Figure 9 for our estimator under the two weighting schemes. Recall that the quadratically decaying weights (6) act essentially like the $\ell_1$ penalty. For numerical comparison, we also include the result of the estimator with $\ell_1$ penalty, which is the CSBS (Convex Sparse Cholesky Selection) method proposed in Khare et al. (2016). For both the non-adaptive banding and the nested lasso methods, we found that their implementations fail to work due to the collinearity of the columns of $X$. 

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Figure 9: Prediction error (computed on an independent test set) of the weighted (left), unweighted (middle) and CSCS (right) estimators.

Figure 9 shows that our estimators are effective in improving modeling performance over a diagonal estimator (attained when $\lambda$ is sufficiently large) and strongly outperform the plain MLE (as evidenced by the sharp increase in prediction error as $\lambda \rightarrow 0$). As expected, the weighted estimator performs very similarly to the CSCS estimator, which uses the $\ell_1$ penalty. Both of these perform better than the unweighted one. However, the sparsity pattern obtained by the two penalties are different (as shown in Figure 10).

Figure 10: Estimates of linkage disequilibrium with tuning parameters selected by 1-se rule and their corresponding precision matrix estimations.

In Figure 10 we show the recovered signed support of the weighted, unweighted and CSCS estimators and their corresponding precision matrices. Black, gray, and white stand for positive, negative, and zero entries respectively. Tuning parameters are chosen using the 1-standard-error rule \cite{Hastie:2009}. The $r$-th row of the estimated matrix $\hat{L}$ reveals the number
of neighboring SNPs necessary for reliably predicting the state of the \( r \)-th SNP. Interestingly, we see some evidence of small block-like structures in \( \hat{L} \), consistent with the hotspot model of recombination as previously described. This regression-based perspective to modeling LD may be a useful complement to the more standard approach, which focuses on raw marginal correlations. Finally, the sparsity recovered by the CSCS estimator, which uses the \( \ell_1 \) penalty, is less easily interpretable, since some entries far from the diagonal are non-zero, losing the notion of ‘local’.

### 6.2 An application to phoneme classification

In this section, we develop an application of our method to a classification problem described in \cite{Hastie2009}. The data consists of \( n = 1717 \) short sound recordings of male voices saying two similar sounding vowels ‘aa’ \((n_1 = 695)\) and ‘ao’ \((n_2 = 1022)\). Each observation \((x_i, y_i)\) has a predictor \( x_i \in \mathbb{R}^p \) representing the (log) intensity of the sound across \( p = 256 \) frequencies and a class label \( y_i \in \{1, -1\} \) for the \( i \)-th observation. Since the frequency is naturally ordered, one might expect that using our estimator in this setting is appropriate.

In linear discriminant analysis (LDA), one assumes that \( x_i | y_i = k \sim N_p(\mu^{(k)}, \Sigma) \) for \( k \in \{1, -1\} \). And in quadratic discriminant analysis (QDA), one further assumes that \( x_i | y_i = k \sim N_p(\mu^{(k)}, \Sigma^{(k)}) \). The LDA/QDA classification rules assign an observation \( x \in \mathbb{R}^p \) to class \( k \) that maximizes \( P(y = k | x) \propto P(x | y = k) \delta(y = k) \), where the estimated probability \( P(x | y = k) \) is calculated using maximum likelihood estimates \( \hat{\mu}^{(k)}, \hat{\Sigma}_i \), and \( \hat{\Sigma}^{(k)} \). More precisely, in the ordered case, the resulting class \( k \) maximizes the LDA/QDA scores:

\[
\begin{align*}
\delta_{\text{LDA}}^{(k)}(x) &= x^T \hat{\Omega}^{(k)} \hat{\mu}^{(k)} - \frac{1}{2} \left( \hat{\mu}^{(k)} \right)^T \hat{\Omega}^{(k)} \hat{\mu}^{(k)} + \log \hat{\pi}^{(k)} = (\hat{L} x)^T \hat{\mu}^{(k)} - \frac{1}{2} \left\| \hat{L} \hat{\mu}^{(k)} \right\|^2 + \log \hat{\pi}^{(k)} \tag{25} \\
\delta_{\text{QDA}}^{(k)}(x) &= x^T \hat{\Omega}^{(k)} \hat{\mu}^{(k)} - \frac{1}{2} \left( \hat{\mu}^{(k)} \right)^T \hat{\Omega}^{(k)} \hat{\mu}^{(k)} + \log \hat{\pi}^{(k)} = (\hat{L}^{(k)} x)^T \hat{\mu}^{(k)} - \frac{1}{2} \left\| \hat{L}^{(k)} \hat{\mu}^{(k)} \right\|^2 + \log \hat{\pi}^{(k)} \tag{26}
\end{align*}
\]

Note that it is the precision matrix, not the covariance matrix, that is used in the above scores. In the setting where \( p > n \), the MLE of \( \Omega \) or \( \Omega^{(k)} \) does not exist. A regularized estimate of precision matrix that exploits the natural ordering information can be helpful in this setting.

To demonstrate the use of our estimator in the high-dimensional setting, we randomly split the data into two parts, with 10\% of the data assigned to the training set and the remaining 90\% of the data assigned to the test set. On the training set, we use 5-fold cross-validation to select the tuning parameter minimizing misclassification error on the validation data. The estimates \( \hat{L} \) and \( \hat{L}^{(k)} \) are then plugged into \( (25) \) and \( (26) \) along with \( \hat{\mu}^{(k)} = \sum_{i \in \text{class } k} x_i / n_k \) and \( \hat{\pi}^{(k)} = n^{(k)} / n_{\text{train}} \) to calculate the misclassification error in the test set. For comparison, we also include non-adaptive banding, the nested lasso, and CSCS. We compute the results (shown in Figure 11) over 10 random train-test splits.

We first observe that, in general, the adaptive methods perform better than the non-adaptive one (which assumes a fixed bandwidth). It is again found that the performance of the weighted estimator is very similar to the one using \( \ell_1 \) penalty (i.e., the CSCS method). And our results are comparable to the nested lasso both in LDA and QDA. Interestingly, we find that the weighted estimator does better in LDA while the unweighted estimator performs better in QDA. The reason, we suspect, is that QDA requires the estimation of more parameters than LDA and therefore favors more constrained methods like the unweighted estimator, which more strongly discourages nonzeros from being far from the diagonal than the weighted one.
An R (R Core Team 2016) package, named varband, will be made available, implementing our estimator. The estimation is very fast with core functions coded in C++, allowing us to solve large-scale problems in substantially less time than is possible with the R-based implementation of the nested lasso.

7 Conclusion

We have presented a new and flexible method of learning local dependence in the setting where the elements of a random vector have a known ordering. The model amounts to sparse estimation of the inverse of the Cholesky factor of the covariance matrix with variable bandwidth. Compared to existing banding-based estimators of the precision matrix in the ordered setting, our method is based on a convex formulation that allows it to simultaneously yield a flexible adaptively-banded sparsity pattern, enjoy efficient computational algorithms, and be studied theoretically. We show how the matrix estimation problem can be decomposed into independent row estimation problems, each of which can be solved via an ADMM algorithm having efficient updates. We prove that our method recovers the signed support of the true Cholesky factor and attains estimation consistency rates in several matrix norms under assumptions as mild as those in linear regression problems. Simulation studies show that our method compares favorably to two pre-existing estimators in the ordered setting, both in terms of support recovery and in terms of estimation accuracy. Through a genetic data example, we illustrate how our method may be applied to model the local dependence of genetic variations in genes along a chromosome. Finally, we illustrate that our method has favorable performance in a sound recording classification problem.

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A Decoupling property

Let \( S = \frac{1}{n} X^T X \in \mathbb{R}^{p \times p} \) be the sample covariance matrix. Then the estimator (4) is the solution to the following minimization problem

\[
\min_{L: L_{rr} > 0 \text{ for } r < k} \left\{ -2 \sum_{r=1}^{p} \log L_{rr} + \text{tr}(SL^T L) + \lambda \sum_{r=2}^{p} \frac{1}{2} \left( \sum_{m=1}^{\ell} w_{rm}^2 L_{rm}^2 \right) \right\}.
\]

First note that under lower-triangular constraint

\[
\text{tr} \left( SL^T L \right) = \frac{1}{n} \sum_{r=1}^{p} \text{tr} \left( XL_{1r}^T L_{rr} X^T \right) = \frac{1}{n} \sum_{r=1}^{p} \| XL_{1r}^T \|_2^2 = \frac{1}{n} \sum_{r=1}^{p} \| X_{1r} L_{1rr}^T \|_2^2,
\]

where \( X_{1r} \) is the matrix of first \( r \) columns of \( X \). Thus

\[
-2 \sum_{r=1}^{p} \log L_{rr} + \text{tr}(SL^T L) + \lambda \sum_{r=2}^{p} \frac{1}{2} \left( \sum_{m=1}^{\ell} w_{rm}^2 L_{rm}^2 \right)
= -2 \log L_{11} + \frac{1}{n} \| X_{11} \|_2^2 + \sum_{r=2}^{p} \left( -2 \log L_{rr} + \frac{1}{n} \| X_{1r} L_{1rr}^T \|_2^2 + \lambda \sum_{m=1}^{\ell} w_{rm}^2 L_{rm}^2 \right).
\]

Therefore the original problem can be decoupled into \( p \) separate problems. In particular, a solution \( \hat{L} \) to can be written in a row-wise form with

\[
\hat{L}_{11} = \arg \min_{L_{11} > 0} \left\{ -2 \log L_{11} + \frac{1}{n} \| X_{11} \|_2^2 \right\} = \frac{1}{\sqrt{S_{11}}},
\]

and for \( r = 2, \ldots, p, \)

\[
\hat{L}_{1rr} = \arg \min_{\beta \in \mathbb{R}^r: \beta_r > 0} \left\{ -2 \log \beta_r + \frac{1}{n} \| X_{1r} \beta \|_2^2 + \lambda \sum_{m=1}^{\ell} w_{rm}^2 \beta_m^2 \right\}.
\]

B A closed-form solution to (8) in the paper

First note that the objective function in (8) in the paper is a smooth function. Take derivative with respect to \( \beta \) and set to zero gives the following system of equations

\[
-2 \frac{1}{\beta_r} e_r + \frac{2}{n} X_{1r}^T X_{1r} \beta + u^{(t-1)} + \rho \left( \beta - \gamma^{(t-1)} \right) = 0.
\]

Denote \( S^{(r)} = \frac{1}{n} X_{1r}^T X_{1r} \), then the equations above can be further decomposed into

\[
-\frac{2}{\beta_r} + \left( 2S^{(r)} + \rho \right) \beta_r + 2S^{(r)} \beta_r + u^{(t-1)} - \rho \gamma^{(t-1)} = 0,
\]

\[
\left( 2S^{(r)} + \rho I \right) \beta_r + 2S^{(r)} \beta_r + u^{(t-1)} - \rho \gamma^{(t-1)} = 0.
\]

Solve for \( \beta_r \) in the second system of equations gives

\[
\beta_r = - \left( 2S^{(r)} + \rho I \right)^{-1} \left( 2S^{(r)} \beta_r + u^{(t-1)} - \rho \gamma^{(t-1)} \right),
\]
where
\[
A = 4S_{r,r}^{(r)} \left( 2S_{r,r}^{(r)} + \rho I \right)^{-1} S_{r,r}^{(r)} - 2S_{r,r}^{(r)} - \rho,
\]
\[
B = 2S_{r,r}^{(r)} \left( 2S_{r,r}^{(r)} + \rho I \right)^{-1} \left( \gamma_{r}^{(t-1)} - \rho \gamma_{r}^{(t-1)} \right) - u_{r}^{(t-1)} + \rho \gamma_{r}^{(t-1)}.
\]

Solving for $\beta_r$ gives the closed update solution.

**C Dual problem of \(\text{(9)}\) in the paper**

**Lemma 7.** A dual problem of \(\text{(9)}\) in the paper is

\[
\min_{\gamma \in \mathbb{R}^r} \left\{ \left\| y^{(t)} - \lambda \sum_{t=1}^{r-1} W(t) * a^{(t)} \right\|_2^2 \right\} \text{ s.t. } \left\| (a^{(t)})_{g_r,t} \right\|_2 \leq 1, \quad (a^{(t)})_{g_r,t} = 0
\]

where $y^{(t)} = \beta^{(t)} + \frac{1}{\rho} u^{(t-1)}$. Also, given a solution $\hat{a}^{(t)}, \ldots, \hat{a}^{(r-1)}$, the solution to \(\text{(9)}\) can be written as

\[
\gamma^{(t)} = y^{(t)} - \lambda \sum_{t=1}^{r-1} W(t) * \hat{a}^{(t)}.
\]

**Proof.** Note that

\[
\sqrt{\sum_{m=1}^{\ell} w_m^2 \gamma_m^2} = \left\| (W^{(t)} * \gamma)_{g_r,t} \right\|_2 = \max \left\{ \left\langle W^{(t)} * a^{(t)}, \gamma \right\rangle, \text{ s.t. } \left\| (a^{(t)})_{g_r,t} \right\|_2 \leq 1, \quad (a^{(t)})_{g_r,t} = 0 \right\}.
\]

So the minimization problem in \(\text{(9)}\) in the paper becomes

\[
\min_{\gamma} \left\{ \frac{1}{2} \left\| \gamma - y^{(t)} \right\|_2^2 + \frac{1}{\rho} \sum_{t=1}^{r-1} \left\| (W(t) * \gamma)_{g_r,t} \right\|_2 \right\} \]

\[
= \max_{a^{(t)}} \left\{ \frac{1}{2} \left\| \gamma - y^{(t)} \right\|_2^2 + \frac{1}{\rho} \sum_{t=1}^{r-1} \left\langle W^{(t)} * a^{(t)}, \gamma \right\rangle, \text{ s.t. } \left\| (a^{(t)})_{g_r,t} \right\|_2 \leq 1, \quad (a^{(t)})_{g_r,t} = 0 \right\}
\]

\[
= \max_{a^{(t)}} \left\{ \min_{\gamma} \left\{ \frac{1}{2} \left\| \gamma - y^{(t)} \right\|_2^2 + \frac{1}{\rho} \sum_{t=1}^{r-1} \left\langle W^{(t)} * a^{(t)}, \gamma \right\rangle, \text{ s.t. } \left\| (a^{(t)})_{g_r,t} \right\|_2 \leq 1, \quad (a^{(t)})_{g_r,t} = 0 \right\} \right\},
\]

where $y^{(t)} = \beta^{(t)} + \frac{1}{\rho} u^{(t-1)}$. Solve the inner minimization problem by setting derivative to zero

\[
\gamma - y^{(t)} + \frac{1}{\rho} \sum_{t=1}^{r-1} W(t) * a^{(t)} = 0,
\]

which gives the primal-dual relation

\[
\gamma = -\frac{1}{\rho} \sum_{t=1}^{r-1} W(t) * a^{(t)} + y^{(t)}.
\]
Plug this back we have

\[
\begin{align*}
\min_\gamma & \left\{ \frac{1}{2} \left\| \gamma - y^{(t)} \right\|^2 + \frac{\lambda}{\rho} \sum_{\ell=1}^{r-1} \left\| \left( W^{(t)} \ast \gamma \right)_{g_{r,\ell}} \right\|_2 \right\} \\
= & \max_{\hat{a}^{(t)}} \left\{ \frac{1}{2} \left\| \frac{-\lambda}{\rho} \sum_{\ell=1}^{r-1} W^{(t)} \ast \hat{a}^{(t)} \right\|_2^2 + \frac{\lambda}{\rho} \sum_{\ell=1}^{r-1} \left( W^{(t)} \ast \hat{a}^{(t)} \right)_{g_{r,\ell}} - \frac{\lambda}{\rho} \sum_{\ell=1}^{r-1} W^{(t)} \ast \hat{a}^{(t)} + y^{(t)} \right\} \\
& \text{s.t.} \left\| \left( \hat{a}^{(t)} \right)_{g_{r,\ell}} \right\|_2 \leq 1, \left( \hat{a}^{(t)} \right)_{g_{r,\ell}} = 0 \right\} \\
= & \min_{\hat{a}^{(t)}} \left\{ \left\| y^{(t)} - \frac{\lambda}{\rho} \sum_{\ell=1}^{r-1} W^{(t)} \ast \hat{a}^{(t)} \right\|_2^2 \right\} \text{s.t.} \left\| \left( \hat{a}^{(t)} \right)_{g_{r,\ell}} \right\|_2 \leq 1, \left( \hat{a}^{(t)} \right)_{g_{r,\ell}} = 0 \right\}.
\end{align*}
\]

\[\square\]

**Algorithm 3** BCD on the dual problem (27)

1. Let $y^{(t)} = \beta^{(t)} + \frac{1}{\rho} \nu^{(t-1)}$
2. Initialize $\hat{a}^{(t)} \leftarrow 0$ for all $\ell = 1, \ldots, r - 1$
3. for $\ell = 1, \ldots, r - 1$ do
4. \indent $\hat{z}^{(t)} \leftarrow y^{(t)} - \frac{\lambda}{\rho} \sum_{k=1}^{r-1} W^{(k)} \ast \hat{a}^{(k)}$ Find a root $\hat{\nu}_\ell$ that satisfies
5. \indent $h_\ell(\nu) := \sum_{m=1}^{\ell} \frac{w_{lm}^2}{(w_{lm}^2 + \nu_\ell)} \left( \hat{z}_m^{(t)} \right)^2 = \frac{\lambda^2}{\rho^2}$ (29)
6. \indent $\hat{a}_m^{(t)} \leftarrow \frac{w_{lm}}{2 \left( w_{lm}^2 + \nu_\ell \right)} \hat{z}_m^{(t)}$
7. \indent return $\{\hat{a}^{(t)}\}$ as a solution to (27)
8. \indent return $\gamma^{(t)} = y^{(t)} - \frac{\lambda}{\rho} \sum_{\ell=1}^{r-1} W^{(t)} \ast \hat{a}^{(t)}$ as a solution to (9) in the paper

**D** Elliptical Projection

We adapt the same procedure as in Appendix B of [Bien et al. (2015)] to update one $a^{(t)}$ in Algorithm (3). By (9) in the paper we need to solve a problem of the form

\[
\min_{\hat{a} \in \mathbb{R}^d} \left\| \hat{z}^{(t)} - \tau D \hat{a} \right\|_2^2 \text{ s.t.} \left\| \hat{a} \right\|_2 \leq 1,
\]

where $\tau = \frac{\lambda}{\rho}$ and $D = \text{diag}(w_{lm})_{m \leq \ell} \in \mathbb{R}^{d \times \ell}$. If $\|D^{-1} \hat{z}^{(t)}\|_2 \leq \tau$, then we could just let $\hat{a} = \frac{1}{\tau} D^{-1} \hat{z}^{(t)}$. Otherwise, we use the Lagrangian multiplier method to solve the constrained minimization problem above. Specifically, we find a stationary point of

\[
\mathcal{L}(a, \nu) = \left\| \hat{z}^{(t)} - \tau D \hat{a} \right\|_2^2 + \nu \tau^2 \left( \left\| \hat{a} \right\|_2^2 - 1 \right).
\]

Taking the derivative with respect to $a$ and set it equal to zero we have

\[
\hat{a}_m = \frac{w_{lm}}{\tau \left( w_{lm}^2 + \nu \right)} \hat{z}_m^{(t)},
\]
for each \(m \leq \ell\) and \(\nu\) is such that \(\|\hat{a}\|_2 = 1\), which means it satisfies (29). By observing that \(h(\nu)\) is a decreasing function of \(\nu\) and \(w_{\ell t} = \max_{m \leq \ell} w_{\ell m}\), following Appendix B of [Bien et al. (2015)](Bien2015), we obtain lower and upper bounds for \(\nu\):

\[
\frac{1}{r} \|D \hat{z}^{(t)}\|_2 - w_{\ell t}^2 \leq \nu \leq \frac{1}{r} \|D \hat{z}^{(t)}\|_2,
\]

which can be used an initial interval for finding \(\nu\) using Newton’s method. In practice, we usually find \(\nu\) from the equation \(\frac{1}{h(\nu)} = \tau^{-2}\) for better numerical stability.

We end this section with a characterization of solution to (9) in the paper, which says that the solution can be written as \(\gamma^{(t)} = y^{(t)} \ast \hat{t}\), where \(\hat{t}\) is some data-dependent vector in \(\mathbb{R}^r\).

**Theorem 8.** A solution to (9) in the paper can be written as \(\gamma^{(t)} = y^{(t)} \ast \hat{g}\), where the data-dependent vector \(\hat{g} \in \mathbb{R}^r\) is given by

\[
\hat{g}_m = \prod_{\ell=m}^{r-1} \frac{[\hat{\nu}_\ell]_+}{w_{\ell m}^2 + [\hat{\nu}_\ell]_+}
\]

and \(\hat{g}_r = 1\), where \(\hat{\nu}_\ell\) satisfies \(\tau^2 = \sum_{m=1}^\ell \frac{w_{\ell m}^2}{(w_{\ell m}^2 + \nu)^2} \left(\hat{z}^{(t)}_{m}\right)^2\).

**Proof.** By [Jenatton et al. (2011)](Jenatton2011), we can get a solution to (9) in the paper in a single pass as described in Algorithm 3. If we start from \(\hat{z}^{(t)} = y^{(t)}\), then for \(\ell = 1, \ldots, r-1\) and each \(m \leq \ell\),

\[
\hat{z}^{(t+1)}_m = \hat{z}^{(t)}_m - \tau w_{\ell m} \hat{a}^{(t)}_m = \frac{[\hat{\nu}_\ell]_+}{w_{\ell m}^2 + [\hat{\nu}_\ell]_+} \hat{z}^{(t)}_m.
\]

By (28), \(\gamma^{(t)} = \hat{z}^{(r-1)}\) and the result follows.

A key observation from this characterization is that a banded sparsity is induced in a solution in (9) in the paper, and thus a solution in (7) in the paper by Algorithm 1.

**Corollary 9.** A solution \(\gamma^{(t)}\) to (9) in the paper has banded sparsity, i.e., \((\gamma^{(t)})_{1:j} = 0\) for \(j = \max \{\ell : \hat{\nu}_\ell \leq 0\}\).

### E Uniqueness of sparse row estimator

**Lemma 10.** (Optimality condition) For any \(\lambda > 0\) and a \(n\)-by-\(p\) sample matrix \(X\), \(\hat{\beta}\) is a solution to the problem

\[
\min_{\beta \in \mathbb{R}^p} \left\{ -2 \log \beta_r + \frac{1}{n} \|X_{1:r}\beta\|_2^2 + \lambda \sum_{\ell=1}^{r-1} \sqrt{\sum_{m=1}^{\ell} w_{\ell m}^2 \beta_m^2} \right\}
\]

iff there exist \(\hat{a}^{(t)} \in \mathbb{R}^r\) for \(\ell = 1, \ldots, r-1\) such that

\[
-\frac{2}{\beta_r} e_r + \frac{2}{n} X_{1:r}^T X_{1:r} \hat{\beta} + \lambda \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \hat{a}^{(t)} = 0
\]

with \((\hat{a}^{(t)})_{g_r,t} = 0\), \((\hat{a}^{(t)})_{g_r,t} = \frac{(W^{(\ell)} \ast \hat{\beta})_{g_r,t}}{\|W^{(\ell)} \ast \hat{\beta}\|_2}\) for \(\hat{\beta}_{g_r,t} \neq 0\) and \(\|\hat{a}^{(t)}\|_2 \leq 1\) for \(\hat{\beta}_{g_r,t} = 0\).
Lemma 11. Take \( \hat{\beta} \) and \( \hat{a}(\ell) \) as in the previous lemma. Suppose that
\[
\left\| \left( \hat{a}(\ell) \right)_{g_r} \right\|_2 < 1 \quad \text{for} \quad \ell = 1, \ldots, J(\hat{\beta})
\]
then for any other solution \( \hat{\beta} \) to \( \hat{\beta} \), it is as sparse as \( \hat{\beta} \) if not more. In other words,
\[
K(\hat{\beta}) \leq \hat{K}_r
\]

Lemma 12. (Uniqueness) Under the conditions of the previous lemma, let \( \hat{S} = \{ i : \hat{\beta}_i \neq 0 \} \). If \( X_S \) has full column rank (i.e., \( \text{rank} (X_S) = |\hat{S}| \)) then \( \hat{\beta} \) is unique.

Proof. See Appendix J, K and L.

\(\square\)

F Proof of Theorem 1

We start with introducing notation. From now on we suppress the dependence on \( r \) in notation for simplicity. We denote the group structure \( g_\ell = \{1, \cdots, \ell\} \) for \( \ell \leq r \) for each \( \ell = 1, \ldots, p \). For any vector \( \beta \in \mathbb{R}^r \), we let \( \beta_{g_\ell} \in \mathbb{R}^\ell \) to be the vector with elements \( \{ \beta_m : m \leq \ell \} \). We also introduce the weight vector \( W(\ell) \in \mathbb{R}^p \) with \( (W(\ell))_m = w_{\ell m} \) where \( w_{\ell m} \) can be defined as in \( \hat{R} \) in the paper or \( w_{\ell m} = 1 \). Finally recall from Section 4 in the paper the definition of \( I \), we denote \( S = I \cup \{ r \} = \{ J + 1, \ldots, r \} \) and \( S^c = \{ 1, 2, \ldots, J \} \).

The general idea of the proof depends on the primal-dual witness procedure in [Wainwright 2009] and [Ravikumar et al. 2011]. Consider the original problem \( \hat{R} \) for any \( r = 2, \ldots, p \), we construct the primal-dual witness solution pairs \( \left( \hat{\beta}, \sum_{\ell=1}^{r-1} W(\ell) \ast \hat{a}(\ell) \right) \) as follows:

(a) Solve the restricted subproblem with the true bandwidth \( K = r - 1 - J \):
\[
\hat{\beta} = \arg \min_{\hat{\beta} \text{ s.t. } \hat{\beta} > 0} \left\{ -2 \log \hat{\beta}_r + \frac{1}{n} \left\| X_{1:r} \hat{\beta} \right\|_2^2 + \lambda \sum_{\ell=1}^{r-1} \left\| (W(\ell) \ast \hat{\beta})_{g_\ell} \right\|_2 \right\}.
\]

Note that without adding any restrictions, the solution above can be written as
\[
\hat{\beta} = \left( 0_{J} \hat{\gamma} \right),
\]
where
\[
\hat{\gamma} = \arg \min_{\gamma \in \mathbb{R}^{K+1}} \left\{ -2 \log \gamma_{K+1} + \frac{1}{n} \left\| X_S \gamma \right\|_2^2 + \lambda \sum_{\ell=1}^{K} \left\| (\hat{W}(\ell) \ast \gamma)_{g_\ell} \right\|_2 \right\},
\]
with
\[
\hat{W}(\ell) = \left( W(\ell+J) \right)_S \iff \sum_{\ell=1}^{K} \left\| (\hat{W}(\ell) \ast \gamma)_{g_\ell} \right\|_2 = \sum_{\ell=J+1}^{r-1} \left( \sum_{m=J+1}^{r-1} w_{\ell m}^2 \right)^{1/2}.
\]

(b) By Lemma 10 there exist \( \check{b}(\ell) \in \mathbb{R}^{K+1} \) for \( \ell = 1, \ldots, K \), such that \( \left( \check{b}(\ell) \right)_{g_\ell} = 0 \) and
\[
\left( \check{b}(\ell) \right)_{g_r} = \frac{\left( \hat{W}(\ell) \ast \hat{\gamma} \right)_{g_\ell}}{\left\| \hat{W}(\ell) \ast \hat{\gamma} \right\|_{g_\ell}}.
\]

satisfying
\[
- \frac{2}{\gamma_{K+1}} e_{K+1} + \frac{2}{n} X^T S \hat{\gamma} + \lambda \sum_{\ell=1}^{K} \hat{W}(\ell) \ast \check{b}(\ell) = 0.
\]
(c) For $\ell = J + 1, \ldots, r - 1$, we let
\[
\hat{a}^{(\ell)} = \begin{pmatrix}
0 \\
\beta^{(\ell-J)}
\end{pmatrix}.
\]

Then we have $(\hat{a}^{(\ell)})_{g_l^\ell} = 0$, \[\frac{\lVert (\hat{a}^{(\ell)})_{g_l^\ell} \rVert}{2} \leq 1, \quad (\hat{a}^{(\ell)})_{g_l^\ell} = \frac{(W^{(\ell)} \hat{\beta})_{g_l^\ell}}{\lVert (W^{(\ell)} \hat{\beta})_{g_l^\ell} \rVert} \quad \text{for } \beta_{g_l^\ell} \neq 0.
\]

(d) Now for each $\ell = 1, \ldots, J$ we choose $\hat{a}^{(\ell)} \in \mathbb{R}^r$ satisfying
\[
\left( \hat{a}^{(\ell)} \right)_\ell = 0 \quad \text{for any } \ell' \neq \ell \quad \text{and} \quad \left( \hat{a}^{(\ell)} \right)_\ell = -\frac{2}{\lambda w_{\ell\ell}} \left( S \hat{\beta} \right)_\ell = -\frac{2}{n\lambda} X_T X S \beta.
\]

Then by construction and the fact that $w_{\ell\ell} = 1$,
\[
\lambda \left( W^{(\ell)} * \hat{a}^{(\ell)} \right)_\ell = \lambda w_{\ell\ell} \left( \hat{a}^{(\ell)} \right)_\ell = -2 \left( S \hat{\beta} \right)_\ell.
\]

Combining with construction in (b)(c) and Lemma 10, \{ $\hat{a}^{(\ell)}$ \} satisfy optimality condition (30):
\[
-\frac{2}{\beta_r} e_r + \frac{2}{n} X^T_1 x_{1:r} \hat{\beta} + \lambda \sum_{\ell=1}^{r-1} W^{(\ell)} * \hat{a}^{(\ell)} = 0
\] (31)

(e) Verify the strict dual feasibility condition for $\ell = 1, \ldots, J$
\[
\left\lVert \frac{2}{n\lambda} X^T_\ell X S \beta \right\rVert = \left\lVert \left( \hat{a}^{(\ell)} \right)_\ell \right\rVert = \frac{\lVert \left( \hat{a}^{(\ell)} \right)_{g_l^\ell} \rVert}{2} < 1.
\] (32)

At a higher level, steps (a) through (d) construct a pair $\left( \hat{\beta}, \{ \hat{a}^{(\ell)} \} \right)$ that satisfy the optimality condition (30), but the \{ $\hat{a}^{(\ell)}$ \} is not necessarily guaranteed to be a member of $\partial \left( P(\hat{\beta}) \right)$. The step (e) verifies the necessary conditions for it to belong to $\partial \left( P(\hat{\beta}) \right)$. Note that by construction in step (b), \{ $\hat{a}^{(\ell)}$ \} satisfies these conditions for $\ell = J + 1, \ldots, r - 1$ since \{ $\hat{b}^{(\ell)}$ \} does, so it remains to verify for $\ell = 1, \ldots, J$ (see step (c)).

For each $\ell = 1, \ldots, J$, by construction in step (d), $(\hat{a}^{(\ell)})_{g_l^\ell} = 0$. Note that $\hat{\beta}_{g^\ell} = 0$ since $\beta_{g^\ell}$ does, for $\hat{a}^{(\ell)}$ to satisfy conditions in Lemma 10 it suffices to show (32).

If the primal-dual witness procedure succeeds, then by construction, the solution $\hat{\beta}$, whose support is contained in the support of true $L_r$, is a solution to (7). Moreover, by strict dual feasibility and Lemma 12, we know that the $\hat{\beta}$ is the unique solution $\hat{\beta}$ to the unconstrained problem (7). Therefore, the support of $\beta$ is contained in the support of $L_r$.

In the following we adapt the same proof technique as [Wainwright (2009)] to show that the primal-dual witness succeeds with high probability, from which we first conclude that $K(\hat{\beta}) \leq K$.

### F.1 Proof of property 1 in Theorem 1

**Proof.** We need to verify the strict dual feasibility (32). By (31),

\[
-\frac{2}{\beta_r} + \frac{2}{n} X^T r X_r \hat{\beta} + \frac{2}{n} X^T r X_r \hat{\beta} = 0,
\]

(i)

\[
\frac{2}{n} X^T r X_r \hat{\beta} + \frac{2}{n} X^T r X_r \hat{\beta} + \lambda \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \hat{a}^{(\ell)} \right)_I = 0.
\]

(ii)
From (ii)

\[ \tilde{\beta}_r = - (X^T I X)^{-1} \left[ X^T I \beta_r + \frac{\lambda_n}{2} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) I \right]. \quad (iii) \]

Plug (iii) back to (i) and denote \( C_I = X^T (X^T X)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) I \) and \( O_I = I - X^T (X^T X)^{-1} X^T \) as the orthogonal projection matrix onto the ortho-complement of column space of \( X_I \), we have

\[ - \frac{2}{\beta_r} + \frac{2}{n} X^T I O_I X \tilde{\beta}_r - \lambda X^T I C_I = 0, \]

which implies that

\[ \tilde{\beta}_r = \frac{\frac{2}{n} X^T I C_I + \sqrt{\frac{\lambda_n}{4} (X^T I C_I)^2 + \frac{4}{n} X^T I O_I X_r}}{\frac{2}{n} X^T I O_I X_r}, \quad (33) \]

and

\[ \left( \tilde{a}^{(\ell)} \right)_\ell = - \frac{2}{n\lambda} X^T I X \tilde{\beta}_S = - \frac{2}{n\lambda} X^T I \tilde{\beta} - \frac{2}{n\lambda} X^T I X \tilde{\beta}_I \]

\[ = - \frac{2}{n\lambda} X^T I \tilde{\beta} + \frac{2}{n\lambda} X^T I (X^T I X)^{-1} \left[ X^T I \tilde{\beta} + \frac{\lambda_n}{2} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) I \right] \]

\[ = - \frac{2}{n\lambda} X^T I \left[ \left( I - X^T (X^T X)^{-1} X^T \right) X \tilde{\beta} + X^T (X^T X)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) \right] \]

\[ = X^T \left[ C_I - O_I \left( \frac{2}{n\lambda} X \tilde{\beta}_r \right) \right]. \quad (34) \]

Conditioning on \( X_I \), we can decompose \( X_r \) and \( X_\ell \) into linear prediction plus prediction error as

\[ X^T I = \Sigma_{rI} (\Sigma_{rI})^{-1} X^T I_\ell + E^T I, \quad (35) \]

\[ X^T I = \Sigma_{rI} (\Sigma_{rI})^{-1} X^T I_\ell + E^T I, \]

where the elements of the error vector \( E_r \sim N \left( 0_n, \theta^{(r)} I_{n \times n} \right) \) and \( E_\ell \sim N \left( 0_n, \theta_\ell^{(r)} I_{n \times n} \right) \), where \( \theta_r^{(\ell)} \) and \( \theta_r^{(r)} \) are defined in Section 4. Then

\[ X^T I O_I = E^T I O_I \quad \text{and} \quad O_I X_r = O_I E_r, \]

and from (34)

\[ \left( \tilde{a}^{(\ell)} \right)_\ell = E^T I \left[ C_I - O_I \left( \frac{2}{n\lambda} E_r \tilde{\beta}_r \right) \right] + \Sigma_{rI} (\Sigma_{rI})^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) \]

\[ := R^{(\ell)} + F^{(\ell)}. \quad (36) \]

We first bound \( \max_{\ell} \left| F^{(\ell)} \right| \). Note that

\[ \left\| \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right\|_I \leq \max_{m \in I} \sum_{\ell=m}^{r-1} \left\| \tilde{a}^{(\ell)} \right\| \leq \max_{m \in I} \sum_{\ell=m}^{r-1} \frac{1}{(\ell - m + 1)^2} = \frac{\pi^2}{6}, \quad (37) \]

\[ \leq \max_{m \in I} \sum_{\ell=m}^{r-1} \left\| \tilde{a}^{(\ell)} \right\| \leq \max_{m \in I} \sum_{\ell=m}^{r-1} \frac{1}{(\ell - m + 1)^2} = \frac{\pi^2}{6}, \]
Lemma 14. For $M$ where the first equality holds from Pythagorean identity. The next lemma bounds the random quantity $a > 0$.

$$\max_{1 \leq i \leq J} \left| \mathbb{E}_{\mathcal{T}} \left( \Sigma_{\mathcal{T}} \right)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right) \right| \leq 1 - \alpha.$$ 

To give a bound on the random quantity $|R^{(\ell)}|$, we first state a general result that will be used multiple times in later proof.

Lemma 13. Consider the term $E_j^T \eta$ where $\eta \in \mathbb{R}^n$ is a random vector depending on $\mathbf{X}_\mathcal{I}$ and $\mathbf{X}_r$ and $E_j \sim \mathcal{N} \left( \mathbf{0}_n, \theta^{(j)} \mathbf{I}_{n \times n} \right)$ for $j = 1, \ldots, J$. If for some $Q \geq 0$

$$P \left[ \text{Var} \left( E_j^T \eta \big| \mathbf{X}_\mathcal{I}, \mathbf{X}_r \right) \geq Q \right] \leq \bar{p}$$

then for any $a > 0$,

$$P \left[ \left| E_j^T \eta \right| \geq a \right] \leq 2 \exp \left( - \frac{a^2}{2Q} \right) + \bar{p}$$

Proof. Define the event

$$\mathcal{B} = \left\{ \text{Var} \left( E_j^T \eta \big| \mathbf{X}_\mathcal{I} \right) \geq Q \right\}.$$

Now for any $a$ and conditioned on $\mathbf{X}_\mathcal{I}$ and $\mathbf{X}_r$,

$$P \left[ E_j^T \eta \geq a \right] \leq P \left[ E_j^T \eta \geq a \big| \mathcal{B}^c \right] + P \left[ \mathcal{B} \right] \leq P \left[ E_j^T \eta \geq a \big| \mathcal{B}^c \right] + \bar{p}.$$

Conditioned on $\mathcal{B}^c$, the variance of $E_j^T \eta$ is at most $Q$. So by standard Gaussian tail bounds, we have

$$P \left[ B^{(r)} \geq a \big| \mathcal{B}^c \right] = E \left[ P \left( B^{(r)} \geq a \big| \mathbf{X}_\mathcal{I}, \mathbf{X}_r \right) \big| \mathcal{B}^c \right] \leq E \left[ 2 \exp \left( - \frac{a^2}{2Q} \right) \big| \mathcal{B}^c \right] \leq 2 \exp \left( - \frac{a^2}{2Q} \right).$$

Then note that $\text{Var} \left( E_{i\ell} \right) = \theta^{(i\ell)} \leq \theta_r$ for $i = 1, \ldots, n$. Now conditioned on both $\mathbf{X}_\mathcal{I}$ and $\mathbf{X}_r$, $R^{(\ell)}$ is zero-mean with variance at most

$$\text{Var} \left( R^{(\ell)} \big| \mathbf{X}_\mathcal{I} \right) \leq \theta_r \left\| \mathbf{C}_\mathcal{I} - \mathbf{O}_\mathcal{I} \left( \frac{2}{n \lambda} E_{r\ell} \tilde{\beta}_r \right) \right\|_2 = \theta_r \left\{ \mathbf{C}_\mathcal{I}^T \mathbf{C}_\mathcal{I} + \left\| \mathbf{O}_\mathcal{I} \left( \frac{2}{n \lambda} E_{r\ell} \tilde{\beta}_r \right) \right\|_2 \right\}$$

$$= \theta_r \left\{ \frac{1}{n} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right)^T \left( \frac{1}{n} \mathbf{X}_\mathcal{I}^T \mathbf{X}_\mathcal{I} \right)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right) _\mathcal{I} + \frac{4}{n^2 \lambda^2} \tilde{\beta}_r^2 \left\| \mathbf{O}_\mathcal{I} E_r \right\|_2 \right\}$$

$$:= \theta_r M_n,$$

where the first equality holds from Pythagorean identity. The next lemma bounds the random quantity $M_n$.

Lemma 14. For $\varepsilon \in (0, \frac{K}{2})$, denote

$$M_n(\varepsilon) := \frac{3\kappa^2 \pi^2 K}{2} \frac{1}{n} + \frac{1}{\theta^{(r)}(n-K)(1-\varepsilon)} + \frac{16}{n^2 \lambda^2},$$

then

$$P \left[ M_n \geq M_n \big| \mathbf{X}_\mathcal{I} \right] \leq 7 \exp \left( -n \min \left\{ \frac{\alpha^2}{4}, \frac{\varepsilon^2}{4} \left( 1 - \frac{K}{n} \right) \right\} \right).$$
Proof. See Appendix [M] \hfill \Box

Now by Lemma [I3] and union bound
\[
P \left[ \max_{1 \leq \ell \leq J} \left| R^{(\ell)} \right| \geq \alpha \right] \leq 2J \exp \left( -\frac{\alpha^2}{2 \theta_r \bar{M}_n(\varepsilon)} \right) + 7 \exp \left( -c_3 n \right), \tag{38}\]
for some constant $c_3$ independent of $n$ and $J$. Note that by assumption $\frac{K}{n} = o(1)$, we have $\bar{M}_n(\varepsilon) \leq 1 - \varepsilon$ for $n$ large enough, thus
\[
\bar{M}_n(\varepsilon) \leq \frac{K}{n} \left( \frac{3n^2 \pi^2}{2} + \frac{1}{K \theta_r(\varepsilon)} (1 - \varepsilon)^2 + 16 K \lambda^2 \right) \leq \frac{K}{n} \left( \frac{3n^2 \pi^2}{2} + \frac{4}{K \theta_r(\varepsilon)} + 16 \right).
\]
For the exponential term in (38), to have faster decaying rate than the $J$ term, we need
\[
\frac{n}{K \log J} > \frac{\theta_r}{\alpha^2} \left( \frac{3n^2 \pi^2}{2} + \frac{8}{K \theta_r(\varepsilon)} + \frac{32}{K \lambda^2} \right).
\]
\hfill \Box

F.2 Proof of property 2 in Theorem [I]

Next we study the $\ell_\infty$ error bound. The following theorem gives $\ell_\infty$ error bound of $\tilde{\beta}$.

Proof. Let $\delta = \tilde{\beta} - \beta^* = \tilde{\beta} - (L^T)_{rr}$, and $W = SL^T - (L)^{-1}$, then from [I] and the fact that $(L)^{-1}$ is lower-triangular,
\[
\delta_{rr} = - (X_r^T X_r)^{-1} \left[ X_r^T X_r \beta_r + (X_r^T X_r) (L^T)_{rr} \right] - \frac{n \lambda}{2} (X_r^T X_r)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right)_{rr} = - \left( \frac{1}{n} X_r^T X_r \right)^{-1} \left[ \frac{1}{n} X_r^T X_r (\delta_r + \beta_r^*) + \frac{1}{n} X_r^T X_r (L^T)_{rr} \right] - \frac{\lambda}{2} \left( \frac{1}{n} X_r^T X_r \right)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right)_{rr}
\]
\[
= (X_r^T X_r)^{-1} X_r^T X_r \delta_r - \frac{1}{n} X_r^T X_r (L^T)_{rr} \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} - \frac{\lambda}{2} \left( \frac{1}{n} X_r^T X_r \right)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \tilde{a}^{(\ell)} \right)_{rr}.
\tag{39}
\]
From [I] and the fact that $(L^{-1})_{rr} = \frac{1}{L_{rr}}$,
\[
- \frac{1}{\beta_r} + \frac{1}{n} X_r^T X_r \delta_r + \frac{1}{X_r^T X_r \delta_r} + \frac{1}{n} X_r^T X_r \beta_r^* + \frac{1}{n} X_r^T X_r \beta_r^* 
\]
\[
= - \frac{1}{\beta_r} + \frac{1}{n} X_r^T X_r \delta_r + \frac{1}{n} X_r^T X_r \delta_r + (SL^T)_{rr} 
\]
\[
= (L^{-1})_{rr} - \frac{1}{\beta_r} + \frac{1}{n} X_r^T X_r \delta_r + \frac{1}{n} X_r^T X_r \delta_r + W_{rr} = 0
\]
\[
= \frac{\delta_r}{L_{rr} \beta_r} + \frac{1}{n} X_r^T X_r \delta_r + \frac{1}{n} X_r^T X_r \delta_r + W_{rr} = 0. \tag{40}
\]
Plugging (39) into (40) we have
\[
\frac{\delta_r}{L_{rr} \beta_r} + \frac{1}{n} X_r^T O \delta_r = X_r^T X_r (X_r^T X_r)^{-1} W_{rr} + \frac{\lambda}{2} X_r C - W_{rr},
\]
\[
\frac{\delta_r}{L_{rr} \beta_r} + \frac{1}{n} X_r^T O \delta_r = X_r^T X_r (X_r^T X_r)^{-1} W_{rr} + \frac{\lambda}{2} X_r C - W_{rr},
\]
Lemma 16. Using the general weighting scheme in the paper, conditioned on \(X_I\), we have

\[
P \left[ \left| X_I^T X_I (X_I^T X_I)^{-1} W_{I,r} \right| \geq \lambda \right] \leq 2 \exp \left( -\frac{2n\lambda^2}{9\theta_1^{(r)} K^2 K} \right) + 2 \exp \left( -\frac{n}{2} \right).
\]

which implies

\[
\delta_r = \left( \frac{1}{L_{rr}} + \frac{1}{n} X_r^T O_r X_r \right)^{-1} \left[ X_r^T X_I (X_I^T X_I)^{-1} W_{I,r} + \frac{1}{2} X_r C_I - W_{rr} \right].
\]

Since \(L_{rr} > 0\) and \(\bar{\beta}_r > 0\),

\[
|\delta_r| \leq \left( \frac{1}{L_{rr}} + \frac{1}{n} X_r^T O_r X_r \right)^{-1} \left( |X_r^T X_I (X_I^T X_I)^{-1} W_{I,r}| + \frac{1}{2} \lambda X_r C_I + |W_{rr}| \right)
\]

Now conditioned on \(X_I\), by the decomposition \([43]\), \((\frac{1}{n} X_r^T O_r X_r)^{-1} = (\frac{1}{n} E_r^T O_I E_r)^{-1} = \frac{1}{\|O_I E_r\|_2^2}\). From lemma \([20]\) it follows that

\[
P \left[ \left( \frac{1}{n} X_r^T O_I X_r \right)^{-1} \geq \frac{1}{\rho^{(r)}} \frac{n}{n-K} \frac{1}{1-\varepsilon} \right] \leq \exp \left( -\frac{1}{4} (n-K)^2 \varepsilon^2 \right).
\]

Also, by Lemma \([19]\)

\[
P \left[ |X_r^T C_I| \geq 1 \right] \leq 2 \exp \left( -\frac{n\alpha^2}{3\theta_1^{(r)} K^2} \right) + 2 \exp \left( -\frac{n}{2} \right).
\]

To deal with the rest of terms in \([40]\) that involve \(W\), we introduce the following concentration inequality to control its element-wise infinity norm.

Lemma 15. Let \(W = S L^T - (L)^{-1}\). Under assumption \([A5]\) and \([A4]\) there exists some constants \(C_1, C_2 > 0\) such that for any \(0 < t \leq 2\kappa\),

\[
P \left[ \|W\|_\infty > t \right] \leq 2p^2 \exp \left( -\frac{C_2 n t^2}{\kappa^2} \right) + 8p \exp \left( -\frac{C_1 n t}{\kappa} \right).
\]

Proof. See Appendix \([N]\) \(\Box\)

Now define the event

\[
A = \{ \|W\|_\infty \leq \lambda \},
\]

then by Lemma \([15]\)

\[
P \left[ A^c \right] \leq 2p^2 \exp \left( -\frac{C_2 n\lambda^2}{\kappa^2} \right) + 8p \exp \left( -\frac{C_1 n \lambda}{\kappa} \right).
\]

The next lemma shows that \(A\) and with the assumption that \(\frac{n^2}{n} = o(1)\), the term \(X_I^T X_I (X_I^T X_I)^{-1} W_{I,r}\) is bounded by a constant times \(\lambda\) with high probability.

Lemma 16. Using the general weighting scheme in the paper, conditioned on \(X_I\), we have

\[
P \left[ \left| X_I^T X_I (X_I^T X_I)^{-1} W_{I,r} \right| \geq \lambda \right] \leq 2 \exp \left( -\frac{2n\lambda^2}{9\theta_1^{(r)} K^2 K} \right) + 2 \exp \left( -\frac{n}{2} \right).
\]
Proof. Recall the decomposition \((35)\)
\[ \mathbf{X}_T^T \mathbf{X}_I (\mathbf{X}_I^T \mathbf{X}_I)^{-1} \mathbf{W}_{I,r} = \Sigma_{rI} (\Sigma_{III})^{-1} \mathbf{W}_{I,r} + E_r^T \mathbf{X}_I (\mathbf{X}_I^T \mathbf{X}_I)^{-1} \mathbf{W}_{I,r}. \]
Conditioned on event \(\mathcal{A}\), by \(\mathcal{A}^3\) and \((37)\)
\[ |\Sigma_{rI} (\Sigma_{III})^{-1} \mathbf{W}_{I,r}| \leq \left\| \Sigma_{rI} (\Sigma_{III})^{-1} \right\| _{\infty} \left\| \mathbf{W}_{I,r} \right\| _{\infty} \leq \lambda. \]
Note that \(\text{Var}(E_{ir}) = \theta_{i}^{(r)}\) for \(i = 1, \ldots, n\). Let \(B^{(r)} := E_r^T \mathbf{X}_I (\mathbf{X}_I^T \mathbf{X}_I)^{-1} \mathbf{W}_{I,r}\), then \(B^{(r)}\) has mean zero and variance at most
\[ \text{Var}(B^{(r)} | \mathbf{X}_I) = \frac{\theta_{i}^{(r)}}{n} \mathbf{W}_r^T \left( \frac{1}{n} \mathbf{X}_I^T \mathbf{X}_I \right)^{-1} \mathbf{W}_r \leq \frac{9\theta_{i}^{(r)} \kappa^2 K \lambda^2}{n}, \]
with probability greater than \(1 - 2 \exp \left( \frac{n}{2} \right)\). Then the result follows from Lemma \([13]\) \(\blacksquare\)

Putting everything together and choosing the tuning parameter from \([12]\) in the paper, with a union bound argument and some algebra, we have shown that conditioned on \(\mathbf{X}_I\),
\[ \mathbf{P} \left[ |\delta_r| \geq \frac{1}{\theta_{i}^{(r)} n - K} \frac{1}{1 - \varepsilon} \frac{5}{2} \lambda \right] \leq \mathbf{P} \left[ |\delta_r| \geq \frac{5}{2\theta_{i}^{(r)}} \lambda \right] = \mathbf{P} \left[ |\delta_r| \geq \frac{5}{2\theta_{i}^{(r)}} \lambda \right] \mathcal{A} + \mathbf{P}[\mathcal{A}^c] \]
\[ \leq \exp \left( -\frac{1}{4n} \left( 1 - \frac{K}{n} \right) \varepsilon^2 \right) + 2 \exp \left( -\frac{\alpha n \theta_{i}^{(r)}}{3\theta_{i}^{(r)} \kappa^2 \pi^2 K} \right) + 2 \exp \left( -\frac{2\alpha n \theta_{i}^{(r)}}{9\theta_{i}^{(r)} \kappa^2 K \lambda^2} \right) + 4 \exp \left( -\frac{n}{2} \right) \]
\[ + 2 P^2 \exp \left( -\frac{C_4 n \lambda^2}{\kappa^2} \right) + 8 \exp \left( -\frac{C_1 n \lambda}{\kappa^2} \right) \leq c_4 \exp \left( -c_3 n \right) + c_6, \quad (41) \]
for some constant \(c_4, c_5, c_6, \varepsilon > 0\) independent of \(n\) and \(p\).

We now consider a bound for \(\delta_{I}\). Recall from \([39]\)
\[ \delta_I = F_1 + F_2 \]
where
\[ F_1 = - (\mathbf{X}_I^T \mathbf{X}_I)^{-1} \mathbf{X}_I^T \mathbf{X}_I \delta_r, \]
\[ F_2 = - \left( \frac{1}{n} \mathbf{X}_I^T \mathbf{X}_I \right)^{-1} \left( \mathbf{W}_{I,r} + \frac{\lambda}{2} \mathbf{D} \right) \quad \text{with} \quad \mathbf{D} = \left( \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \hat{q}^{(\ell)} \right). \]

Since
\[ \|F_2\|_{\infty} \leq \left\| \left( \frac{1}{n} \mathbf{X}_I^T \mathbf{X}_I \right)^{-1} - (\Sigma_{III})^{-1} \right\| _{\infty} \left\| \left( \mathbf{W}_{I,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\| _{\infty} + \left\| (\Sigma_{III})^{-1} \left( \mathbf{W}_{I,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\| _{\infty}, \]
where the second term, by \((37)\)
\[ \left\| (\Sigma_{III})^{-1} \left( \mathbf{W}_{I,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\| _{\infty} \leq \left\| (\Sigma_{III})^{-1} \right\| _{\infty} \left\| \mathbf{W}_{I,r} \right\| _{\infty} + \frac{\lambda}{2} \left\| \mathbf{D} \right\| _{\infty} \]
\[ \leq \left\| (\Sigma_{III})^{-1} \right\| _{\infty} \left( 1 + \frac{\pi^2}{12} \right) \lambda \leq 2 \lambda \left\| (\Sigma_{III})^{-1/2} \right\| _{\infty}^2 \quad \text{on} \ \mathcal{A}. \]

To deal with the first term, note that \(X_I = W_I (\Sigma_{III})^{1/2}\), where \(W_I \in \mathbb{R}^{n \times K}\) is a standard Gaussian random matrix (i.e., \((W^T_I)_{ij} \sim N(0,1))\). Thus we can write it as
\[ \left\| (\Sigma_{III})^{-1/2} \left( \frac{1}{n} W_I^T W_I \right)^{-1} - I_K \right\| _{\infty} \right\| (\Sigma_{III})^{-1/2} \left( \mathbf{W}_{I,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\| _{\infty} \leq \left\| (\Sigma_{III})^{-1/2} \right\| _{\infty} G, \]
where
\[
G = \left\| \frac{1}{n} W_{x}^T W_{x} \right\|^{-1} - \mathbf{I}_{K} \right\|_{\infty} (\Sigma_{II})^{-1/2} \left( W_{x,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\|_{\infty}.
\]

By Lemma 5 in [Wainwright (2009)], we have, for some constant \(c_7 > 0\).
\[
P \left[ G \geq \left\| (\Sigma_{II})^{-1/2} \left( W_{x,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\|_{\infty} \right] \leq 4 \exp \left( -c_7 \min \{ K, \log J \} \right).
\]

Note that conditioning on \(A\), \(\left\| (\Sigma_{II})^{-1/2} \left( W_{x,r} + \frac{\lambda}{2} \mathbf{D} \right) \right\|_{\infty}\) is upper bounded by \(2\lambda\left\| (\Sigma_{II})^{-1/2} \right\|_{\infty} \). Thus
\[
P \left[ G \geq 2\lambda\left\| (\Sigma_{II})^{-1/2} \right\|_{\infty} A \right] \leq 4 \exp \left( -c_7 \min \{ K, \log J \} \right),
\]
and
\[
P \left[ \left\| F_2 \right\|_{\infty} \geq 4\lambda\left\| (\Sigma_{II})^{-1/2} \right\|_{\infty} \right] \leq \frac{P \left[ \left\| F_2 \right\|_{\infty} \geq 4\lambda\left\| (\Sigma_{II})^{-1/2} \right\|_{\infty} \right]}{P \left[ A^c \right]} 
\leq 4 \exp \left( -c_7 \min \{ K, \log J \} \right) + \frac{c_6}{p}.
\] (42)

Turn to \(F_1\), conditioned on \(X_{x}\), by decomposition (35)
\[
\left\| F_1 \right\|_{\infty} \leq \left\| (\Sigma_{II})^{-1} \Sigma_{x} \right\|_{\infty} \left\| |\delta_r| + (X_{x}^T X_{x})^{-1} X_{x}^T E_r| \right\|_{\infty}.
\]

By (41) and [A3]
\[
P \left[ \left\| (\Sigma_{II})^{-1} \Sigma_{x} \right\|_{\infty} \left\| |\delta_r| \geq \frac{5}{2\theta_r^{(r)}} \right\| \leq c_4 \exp \left( -c_5 n \right) + \frac{c_6}{p}.
\]

Consider each coordinate \(j \in I\) of random term whose variance is bounded by
\[
\Var \left[ e_j^T \left( X_{x}^T X_{x} \right)^{-1} X_{x}^T E_r| X_{x} \right] \leq \theta_r \left\| \left( X_{x}^T X_{x} \right)^{-1} \frac{\delta_r}{n} \right\|_{2}.
\]

By Lemma [18] and (41)
\[
P \left[ \left\| \left( X_{x}^T X_{x} \right)^{-1} X_{x}^T E_r \right\|_{\infty} \geq \frac{5}{2\theta_r^{(r)}} \right\| \leq 2 \exp \left( -\frac{n}{4 \theta_r \kappa^2} \right) + c_4 \exp \left( -\frac{c_5 n}{2} \right) + \frac{c_6}{p}.
\]

Thus by Lemma [13]
\[
P \left[ \left\| \left( X_{x}^T X_{x} \right)^{-1} X_{x}^T E_r \right\|_{\infty} \geq \frac{5}{2\theta_r^{(r)}} \right\| \leq 2 \exp \left( -\frac{n}{18 \theta_r \kappa^2} \right) + c_4 \exp \left( -\frac{c_5 n}{2} \right) + \frac{c_6}{p},
\]
and
\[
P \left[ \left\| F_1 \right\|_{\infty} \geq \frac{5}{\theta_r^{(r)}} \right\| \leq 2 \exp \left( -\frac{n}{18 \theta_r \kappa^2} \right) + 2 \exp \left( -\frac{n}{2} \right) + c_4 \exp \left( -\frac{c_5 n}{2} \right) + \frac{c_6}{p}.
\]

Combining with (41) and (42) we have
\[
P \left[ \left\| \delta \right\|_{\infty} \geq 4\lambda\left\| (\Sigma_{II})^{-1/2} \right\|_{\infty} \right] \leq c_8 \exp \left( -c_9 n \right) + \frac{c_6}{p} + 4 \exp \left( -c_7 \min \{ K, \log J \} \right),
\]
for some constants \(c_8, c_9 \geq 0\) independent of \(n\) and \(J\).
Finally we establish a \( \beta_{\min} \) condition, which, combined with \( \ell_\infty \) rate, should give the other direction of the support recovery, i.e., \( K(\hat{\beta}) \geq K \).

By the triangular inequality

\[
|\hat{\beta}_j| \geq |\beta_j| - |\tilde{\beta}_j - \beta_j|.
\]

So if we have

\[
\max_{j \geq J + 1} \left\{ |\beta_j| - |\tilde{\beta}_j - \beta_j| \right\} > 0,
\]

then \( K(\tilde{\beta}) \geq K \).

**G Proof of Theorem 3**

**Proof.** The overall proving techniques are the same as the proof of Theorem 1. The first part of the Theorem holds if \( \max_{2 \leq r \leq p} \max_{1 \leq \ell \leq J_r} |\tilde{a}^{(r\ell)}| < 1\). Now for each \( r = 2, \ldots, p \) we proceed with the same primal-dual witness procedure and end up with the same decomposition (36).

The irrepresentable assumption \( A3 \) ensures that \( \max_{2 \leq r \leq p} \max_{1 \leq \ell \leq J_r} |F^{(r\ell)}| \leq 1 - \alpha \). Following the same line of proof to deal with random term \( R^{(r\ell)} \), we have that \( R^{(r\ell)} \) is zero-mean Gaussian with conditional variance bounded above by the scaling

\[
\theta_r M_n^{(r)}(\varepsilon) = \frac{3\kappa^2 \pi^2 \theta_r K_r^*}{n} + \frac{\theta_r}{\theta_r^{(r)}} \frac{1}{(n - K_r^*) (1 - \varepsilon)} + \frac{16\theta_r}{n\lambda^2} \leq \frac{3\kappa^2 \pi^2 \theta_r}{2} \left( \frac{K}{n} + \frac{\kappa^2}{n\theta_r^{(r)}} (1 - \varepsilon)^2 + \frac{16}{n\lambda^2} \right),
\]

for \( \varepsilon \in (0, \frac{1}{2}) \) with high probability where we use the fact that \( K = o(n) \) thus \( \frac{K}{n} \leq \varepsilon \) for \( n \) large. And

\[
P \left[ \left| R^{(r\ell)} \right| \geq \alpha \right] \leq 2 \exp \left( -\frac{\alpha^2}{2\theta_r M_n^{(r)}(\varepsilon)} \right) + 7 \exp (-c_3 n).
\]

Thus

\[
P \left[ \max_{2 \leq r \leq p} \max_{1 \leq \ell \leq J_r} \left| R^{(r\ell)} \right| \geq \alpha \right] \leq 2 \sum_{r=2}^{p} J_r \exp \left( -\frac{\alpha^2}{2\theta_r M_n^{(r)}(\varepsilon)} \right) + 7 \sum_{r=2}^{p} J_r \exp (-c_3 n) \leq p^2 \exp \left( -\frac{\alpha^2}{3\kappa^2 \pi^2 \theta K n} + \frac{8\theta \kappa^2}{n} + \frac{32\theta \lambda^2}{n^2} \right) + 7 \frac{p^2}{2} \exp (-c_3 n).
\]

For the exponential term to decay faster than \( p^2 \), we need

\[
\frac{n}{\log p} > \max \left\{ \frac{2}{\alpha^2} \left( 3\kappa^2 \pi^2 \theta K + 8\kappa^2 \theta + \frac{32\theta}{\lambda^2} \right), \frac{2}{c_3} \right\}.
\]
H  Proof of Theorem 4

Corollary 17. Using notation and conditions in Theorem 3, the following deviation bounds hold with high probability:

\[ \| \hat{L} - L \|_\infty \leq \zeta \sqrt{\frac{n \log p}{n}}, \]
\[ \| \hat{L} - L \|_1 \leq \zeta (K + 1) \sqrt{\frac{n \log p}{n}}, \]
\[ \| \hat{L} - L \|_2 \leq \zeta (K + 1) \sqrt{\frac{n \log p}{n}}, \]
\[ \| \hat{L} - L \|_F \leq \zeta \sqrt{(s + p) \log p} \frac{n}{n}. \]

Proof. By Theorem 3 with high probability, the support of \( \hat{L} \) is contained in the true support and

\[ \| \hat{L} - L \|_\infty \leq \zeta \sqrt{\frac{n \log p}{n}}. \]

Note that

\[ \| \hat{L} - L \|_\infty = \max_{2 \leq r \leq p} \sum_{c=1}^r | \hat{L}_{rc} - L_{rc} | \leq \max_{1 \leq r \leq p} (K_r + 1) \| \hat{L} - L \|_\infty \leq (K + 1) \| \hat{L} - L \|_\infty. \]

Denote \( D = \max_{1 \leq c \leq p-1} D_c \) where \( D_c = | \{ r = c, \ldots, p : L_{rc} \neq 0 \} | \). Observe that \( D \leq K \), we have

\[ \| \hat{L} - L \|_1 = \max_{1 \leq r \leq p} \sum_{c=1}^r | \hat{L}_{rc} - L_{rc} | \leq \max_{1 \leq r \leq p} (D_c + 1) \| \hat{L} - L \|_\infty \leq (D + 1) \| \hat{L} - L \|_\infty \leq (K + 1) \| \hat{L} - L \|_\infty. \]

By Hölder’s inequality

\[ \| \hat{L} - L \|_2 \leq \sqrt{\| \hat{L} - L \|_1 \| \hat{L} - L \|_\infty}. \]

For Frobenius norm, since

\[ \| \hat{L} - L \|_F^2 = \sum_{r=2}^p \sum_{c=J_r+1}^r (\hat{L}_{rc} - L_{rc})^2 \leq \sum_{r=2}^p \sum_{c=J_r+1}^r \| \hat{L} - L \|_\infty^2 \leq \zeta^2 (\sum_r K_r + p) \frac{n \log p}{n}. \]

of Theorem 4. First note that

\[ \hat{L}^T \hat{L} - L^T L = (\hat{L} - L)^T (\hat{L} - L) + \hat{L}^T L + L^T \hat{L} - 2L^T L \]
\[ = (\hat{L} - L)^T (\hat{L} - L) + (\hat{L} - L)^T L + L^T (\hat{L} - L). \]

Thus,

\[ \| \hat{L}^T \hat{L} - L^T L \|_\infty \leq \| \hat{L} - L \|_\infty \| \hat{L} - L \|_\infty + 2 \| L \|_\infty \| \hat{L} - L \|_\infty, \]
\[ \| \hat{L}^T \hat{L} - L^T L \|_1 \leq \| \hat{L}^T L - L^T L \|_\infty + \| \hat{L} - L \|_\infty^2. \]

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By Hölder’s inequality
\[
\|\hat{L}^T \hat{L} - L^T L\|_2 \leq \sqrt{\|\hat{L}^T \hat{L} - L^T L\|_1 \|\hat{L}^T \hat{L} - L^T L\|_\infty}.
\]

Finally, for Frobenius norm, observe that
\[
\|L^T (\hat{L} - L)\|_F = \|\text{vec} \left( L^T (\hat{L} - L) \right)\|_2 = \|(I_p \otimes L^T) \text{vec} (\hat{L} - L)\|_2
\leq \|I_p \otimes L^T\|_2 \|\hat{L} - L\|_F = \|L\|_2 \|\hat{L} - L\|_F,
\]
where the last equality holds since \(L\) is lower-triangular. Applying the same strategy to
\[
\|L^T (\hat{L} - L) (\hat{L} - L)\|_F,
\]
we have
\[
\|\hat{L}^T \hat{L} - L^T L\|_F \leq \left( \|\hat{L} - L\|_2 + 2\|L\|_2 \right) \|\hat{L} - L\|_F,
\]
then the results follow from Corollary 17.

I Proof of Theorem 6

Proof. Let
\[
G(\Delta) = -2 \log \det (L + \Delta) + \text{tr} \left( S (L + \Delta)^T (L + \Delta) \right) + \lambda \|\Delta + L\|_{2,1}^* + 2 \log \det L - \text{tr} \left( S L^T L \right) - \lambda \|L\|_{2,1}^*,
\]
where \(L\) is the Cholesky factor of the true precision matrix \(\Omega = L^T L\), and the penalty is defined above as
\[
\|L\|_{2,1}^* = \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} \sqrt{\sum_{m=1}^{\ell} w_{lm}^2 L_{rm}^2}.
\]

Note that the estimator \(\hat{L}\) is defined as
\[
\hat{L} = \arg \min_{L_{jk} = 0, j < k} \left\{ -2 \log \det L + \text{tr} \left( S L^T L \right) + \lambda \|L\|_{2,1}^* \right\}.
\]

Thus, \(\hat{\Delta} = \hat{L} - L\) minimize \(G(\Delta)\). Note that we suppress the dependence on \(\lambda\) in \(\hat{L}\) and \(\hat{\Delta}\). We adapt the main idea of the proof in Rothman et al. (2008). Consider the value of \(G(\Delta)\) on the set defined as
\[
\Theta_n(M) = \{\Delta : \Delta_{jk} = 0 \text{ for } j < k, \|\Delta\|_F = Mr_n\},
\]
where \(M > 0\) and
\[
r_n = \sqrt{\frac{\sum_{r=2}^{p} K_r + p}{n}} \log p \rightarrow 0.
\]

Note that
\[
G(\hat{\Delta}) \leq G(0) = 0,
\]
and by the convexity of \(G(\Delta)\), if we can show that
\[
\inf \{G(\Delta) : \Delta \in \Theta_n(M)\} > 0,
\]
then the results follow from Corollary 17.
the minimizer $\hat{\Delta}$ must be inside the sphere defined by $\Theta_n(M)$ and hence
\[
\|\hat{\Delta}\|_F = \|\hat{L} - L\|_F \leq Mr_n.
\]

For the logarithm term in (43), we let $f(t) = \log \det (L + t\Delta)$. Using Taylor expansion of $f(t)$ and the integral form of the remainder with $f'(t) = \text{tr}[(L + t\Delta)^{-1}\Delta]$ and $f''(t) = -\text{vec} \Delta^T (L + t\Delta)^{-1} \otimes (L + t\Delta)^{-1} \text{vec} \Delta$, we have
\[
\log \det (L + \Delta) - \log \det (L) = \text{tr}(L^{-1}\Delta) - (\text{vec} \Delta)^T \left[ \int_0^1 (1 - \nu)(L + \nu \Delta)^{-1} \otimes (L + \nu \Delta)^{-1} d\nu \right] (\text{vec} \Delta).
\]

The trace term in (43) can be written as
\[
\text{tr} \left( S (L + \Delta)^T (L + \Delta) \right) - \text{tr} (SL^T L) = \text{tr} \left( SL^T \Delta + S \Delta^T L + S \Delta^T \Delta \right)
= 2 \text{tr} \left( SL^T \Delta \right) + \text{tr} \left( S \Delta^T \Delta \right)
\geq 2 \text{tr} \left( SL^T \Delta \right),
\]
where the last inequality comes from the fact that the sample covariance matrix $S$ is positive semi-definite. Combining these with (43) gives
\[
G(\Delta) \geq 2(\text{vec} \Delta)^T \left[ \int_0^1 (1 - \nu)(L + \nu \Delta)^{-1} \otimes (L + \nu \Delta)^{-1} d\nu \right] (\text{vec} \Delta)
+ 2 \text{tr}[SL^T - L^{-1}]\Delta + \lambda \left( \|L + \Delta\|_{\Sigma,1}^2 - \|L\|_{\Sigma,1}^2 \right)
\equiv (a) + (b) + (c).
\]

The integral term $(a)$ above has a positive lower bound. Recall that $\sigma_{\text{min}}(M) = \min_{\|x\|_1} x^T M x$ is a concave function of $M$ (the minimum of linear functions of $M$ is concave), we have
\[
(a) = 2\|\text{vec} \Delta\|^2 \text{vec} \Delta^T \left[ \int_0^1 (1 - \nu)(L + \nu \Delta)^{-1} \otimes (L + \nu \Delta)^{-1} d\nu \right] \text{vec} \Delta
\geq 2\|\Delta\|_F^2 \sigma_{\text{min}} \left[ \int_0^1 (1 - \nu)(L + \nu \Delta)^{-1} \otimes (L + \nu \Delta)^{-1} d\nu \right]
\geq 2\|\Delta\|_F^2 \int_0^1 (1 - \nu)\sigma_{\text{min}}^2(L + \nu \Delta)^{-1} d\nu
\geq \|\Delta\|_F^2 \min_{0 \leq \nu \leq 1} \sigma_{\text{min}}^2(L + \nu \Delta)^{-1}
\geq \|\Delta\|_F^2 \min \left\{ \sigma_{\text{min}}^2(L + \Delta)^{-1} : \|\Delta\|_F \leq Mr_n \right\}.
\]

The second inequality uses the Jenson’s Inequality of the concave function $\sigma_{\text{min}}()$ and the fact that the eigenvalues of Kronecker products of lower-triangular matrices are the products of the eigenvalues of their factors. Now
\[
\sigma_{\text{min}}^2(L + \Delta)^{-1} = \sigma_{\text{max}}^2(L + \Delta) \geq (\|L\|_2 + \|\Delta\|_2)^{-2} \geq \frac{1}{2\|L\|_2^2} \geq \frac{\kappa^2}{2},
\]
with probability tending to 1 since $\|L\|_2 \leq \|\Delta\|_F = o(1)$ and $\Lambda 4$. This gives the lower bound for the first term in (45)
\[
(a) \geq \frac{1}{2} \kappa^2 \|\Delta\|_F^2 = \frac{1}{2} \kappa^2 M^2 \kappa_n^2.
\]
To deal with (b), we start by recalling some notation. We denote \(S = \{(r, j) : L_{rj} \neq 0\}\) be the support of \(L\), and \(s = \sum_{r=2}^{p} K_r\) be the number of nonzero off-diagonal elements. We also define
\[
\|L\|_{2,1} = \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} w_{\ell \ell} |L_{r\ell}| = \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} |L_{r\ell}|,
\]
where the last equality holds since \(w_{\ell \ell} = 1\) by (6). Then, by Cauchy-Schwarz inequality,
\[
|\text{tr}((SL^T - L^{-1})\Delta)| = \left| \sum_{r=1}^{p} \sum_{j=1}^{r} (SL^T - L^{-1})_{rj} \Delta_{rj} \right|
\]
\[
\leq \sum_{r=1}^{p} \sum_{j \in I_r} (SL^T - L^{-1})_{rj} \Delta_{rj} + \sum_{r \in I_r} \sum_{j \notin I_r} (SL^T - L^{-1})_{rj} \Delta_{rj}
\]
\[
\leq \sqrt{s + p} \|SL^T - L^{-1}\|_\infty \|\Delta_S\|_F + \|SL^T - L^{-1}\|_\infty \|\Delta_{S^c}\|_{2,1}
\]
\[
\leq C_1 \sqrt{s + p} \left( \frac{\log p}{n} \|\Delta_S\|_F + \frac{\log p}{n} \|\Delta_{S^c}\|_{2,1} \right),
\]
where the last inequality comes from Lemma 15 with probability tending to 1. To bound the penalty terms, we note that
\[
2\lambda \|\Delta_{S^c}\|_{2,1} = \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} \sqrt{\sum_{m:(r,m) \in S} w_{\ell m}^2 (L_{rm} + \Delta_{rm})^2 + \sum_{m:(r,m) \notin S} w_{\ell m}^2 (L_{rm} + \Delta_{rm})^2 - \|L_S\|_{2,1}^2}
\]
\[
\geq \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} \sqrt{\sum_{m:(r,m) \in S} w_{\ell m}^2 (L_{rm} + \Delta_{rm})^2 + \sum_{m:(r,m) \notin S} w_{\ell m}^2 (L_{rm} + \Delta_{rm})^2 - \|L_{r\ell} + \Delta_{r\ell}\| - \|L_S\|_{2,1}^2}
\]
\[
= \|L_S + \Delta_S\|_{2,1}^2 + \|L_{S^c} + \Delta_{S^c}\|_{2,1}^2 - \|L_S\|_{2,1}^2
\]
\[
\geq \|\Delta_{S^c}\|_{2,1}^2 - \|\Delta_{S^c}\|_{2,1}^2
\]
where the last inequality comes from triangle inequality. To give an upper bound on \(\|L_S\|_{2,1}\), by the observation that \(2\lambda b \leq a \lambda^2 + b^2 / a\) holds for any \(a > 0\), we obtain
\[
2\lambda \|\Delta_{S^c}\|_{2,1}^2 = \sum_{r=2}^{p} 2\lambda \sum_{\ell=J_{r+1}}^{r-1} \sqrt{\sum_{m=J_r+1}^{\ell} w_{\ell m}^2 \Delta_{rm}^2}
\]
\[
\leq \left( \sum_{r=2}^{p} K_r \right) \lambda^2 a + \sum_{r=2}^{p} \sum_{\ell=J_{r+1}}^{r-1} \sum_{m=J_r+1}^{\ell} w_{\ell m}^2 \Delta_{rm}^2 / a
\]
\[
= \left( \sum_{r=2}^{p} K_r \right) \lambda^2 a + \sum_{r=2}^{p} \sum_{\ell=J_{r+1}}^{r-1} \left( \sum_{m=J_r+1}^{\ell} w_{\ell m}^2 \right) \Delta_{rm}^2 / a.
\]
Now let
\[
a = \frac{4}{k^2} \max_{r} \max_{J_{r+1} \leq m \leq r-1} \sum_{\ell=m}^{r-1} w_{\ell m}^2
\]
\[
= \frac{4}{k^2} \max_{r} \max_{J_{r+1} \leq m \leq r-1} \sum_{\ell=m}^{r-1} \frac{1}{(\ell - m + 1)^4} \leq \frac{4}{k^4 k^2} \leq \frac{C_2}{k^2},
\]
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for some constant $C_2 > 0$, it follows that
\[
\lambda \|s\|_{2,1}^2 \leq \frac{C_2}{\kappa^2} s^2 \lambda^2 + \|s\|_F^2 \frac{\kappa^2}{4} \leq \frac{C_2}{\kappa^2} s^2 \lambda^2 + \|s\|_F^2 \frac{\kappa^2}{4}.
\]

Therefore,
\[
\lambda \left( \|L + \Delta\|_{2,1}^2 - \|L\|_{2,1}^2 \right) \geq \lambda \|s\|_{2,1}^2 - \frac{C_2}{\kappa^2} s^2 \lambda^2 - \frac{\kappa^2}{4} \|s\|_F^2.
\]
(48)

Finally, combining (46), (47) and (48), we have
\[
G(\Delta) \geq \frac{\kappa^2}{4} \|s\|_F^2 - C_1 \sqrt{\frac{(s + p) \log p}{n}} \|\Delta\|_F + \left( \lambda - C_1 \sqrt{\frac{\log p}{n}} \right) \|s\|_{2,1}^2 - \frac{C_2}{\kappa^2} s^2 \lambda^2.
\]

Now take
\[
\lambda = \frac{C_1}{\varepsilon} \sqrt{\frac{\log p}{n}},
\]
for $\varepsilon < 1$ and recall that $\|\Delta\|_F = Mr_n$, we have
\[
G(\Delta) \geq \frac{\kappa^2}{4} M^2 r_n^2 - C_1 M r_n^2 + C_1 \sqrt{\frac{\log p}{n}} \left( \frac{1}{\varepsilon} - 1 \right) \|s\|_F^2 - \frac{C_2 C^2}{\kappa^2 \varepsilon^2} \frac{s \log p}{n}
\geq \left( \frac{\kappa^2}{4} M^2 - C_1 M - \frac{C_2 C^2}{\kappa^2 \varepsilon^2} \right) r_n^2 > 0,
\]
for $M$ sufficiently large. This establishes the theorem.

\[
J \quad \text{Proof of Lemma 10}
\]

\textbf{Proof.} Denote
\[
\mathcal{L}(\tau, z, \beta; \nu, \phi, a^{(t)}) = -2 \log \tau + \frac{1}{n} \|z\|_2^2 + \nu (\tau - \beta) + \frac{1}{n} \langle \phi, z - X_{1; r, \beta} \rangle + \lambda \sum_{t=1}^{r-1} \langle W^{(t)} * a^{(t)}, \beta \rangle.
\]

Then the primal (7) in the paper can be written equivalently as
\[
\min_{\tau, z, \beta} \left\{ \max_{\nu, \phi, a^{(t)}} \left\{ \mathcal{L}(\tau, z, \beta; \nu, \phi, a^{(t)}): \left\| (a^{(t)})_{g_r, t} \right\|_2 \leq 1, \left( a^{(t)} \right)_{g_r, t} = 0 \right\} \right\}.
\]

Then the dual function can then be written as
\[
g(\nu, \phi, a^{(t)}) = \inf_{\tau, z, \beta} \mathcal{L}(\tau, z, \beta; \nu, \phi, a^{(t)})
= \inf_{\tau} \{-2 \log \tau + \nu \tau\} + \inf_{\varepsilon} \left\{ \frac{1}{n} \|z\|_2^2 + \frac{1}{n} \langle \phi, z \rangle \right\} + \inf_{\beta} \left\{ -\nu \beta + \frac{1}{n} \langle X_{1; r, \beta} \rangle + \lambda \sum_{t=1}^{r-1} \langle W^{(t)} * a^{(t)}, \beta \rangle \right\}
= 2 \log \nu - 2 \log 2 + 2 + \infty \{ \nu > 0 \} - \frac{1}{4n} \|\phi\|_2^2 - \infty \left\{ -\nu e_r - \frac{1}{n} \langle X_{1; r, \phi} \rangle + \lambda \sum_{t=1}^{r-1} \langle W^{(t)} * a^{(t)}, \beta \rangle = 0 \right\},
\]
where $e_r \in \mathbb{R}^r$ such that $(e_r)_r = 1$ and $(e_r)_j = 0$ for all $j \neq r$. Thus the dual problem (by a constant difference) is
\[
\max_{\nu, \phi, a^{(t)}} g(\nu, \phi, a^{(t)})
= \min_{\nu, \phi, a^{(t)}} \left\{ -2 \log \nu - \frac{1}{4n} \|\phi\|_2^2 \quad \text{s.t.} \quad \nu > 0, \quad \left\| (a^{(t)})_{g_r, t} \right\|_2 \leq 1, \left( a^{(t)} \right)_{g_r, t} = 0, \quad \nu e_r + \frac{1}{n} \langle X_{1; r, \phi} \rangle = \lambda \sum_{t=1}^{r-1} \langle W^{(t)} * a^{(t)} \rangle \right\}.
\]

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And the primal-dual relation is
\[ \hat{\beta}_r = \hat{\tau} = \frac{2}{\nu} \quad \hat{\phi} = -2\hat{z} = -2X_{1,r}\hat{\beta}. \]

This implies that at optimal points
\[ -\frac{2}{\hat{\beta}_r}e_r + 2S_{1,r,1:r}\hat{\beta} + \lambda \sum_{\ell=1}^{r-1} W^{(\ell)} \ast \hat{a}^{(\ell)} = 0, \]
with \( \left\| (\hat{a}^{(\ell)})_{g_{r,\ell}} \right\|_2 \leq 1, (\hat{a}^{(\ell)})_{g_{r,\ell}} = 0. \)

If we denote the objective function as
\[ f(\beta) = -2\log\beta_r + \langle S_{1:r,1:r}, \beta\beta^T \rangle + \lambda P(\beta), \]
then from the equality \( f(\hat{\beta}) = \mathcal{L}(\hat{\tau}, \hat{z}, \hat{\beta}; \hat{\nu}, \hat{\phi}, \hat{a}^{(\ell)}) \) together with the primal-dual relation, we have
\[ P(\hat{\beta}) = \sum_{\ell=1}^{r-1} \langle W^{(\ell)} \ast \hat{a}^{(\ell)}, \hat{\beta} \rangle = \sum_{\ell=1}^{r-1} \langle W^{(\ell)} \ast \hat{\beta}, \hat{a}^{(\ell)} \rangle. \]

Suppose there exists some \( \ell \) with \( \hat{\beta}_{g_{r,\ell}} \neq 0 \) but \( (\hat{a}^{(\ell)})_{g_{r,\ell}} \neq \frac{\langle W^{(\ell)} \ast \hat{\beta}, \hat{a}^{(\ell)} \rangle}{\left\| (W^{(\ell)} \ast \hat{\beta})_{g_{r,\ell}} \right\|_2} \), then \( \langle W^{(\ell)} \ast \hat{\beta}, \hat{a}^{(\ell)} \rangle < \left\| (W^{(\ell)} \ast \hat{\beta})_{g_{r,\ell}} \right\|_2 \) while for other \( \ell' \) by Cauchy-Schwarz inequality we have \( \langle W^{(\ell')} \ast \hat{\beta}, \hat{a}^{(\ell')} \rangle \leq \left\| (W^{(\ell')} \ast \hat{\beta})_{g_{r,\ell'}} \right\|_2 \). Therefore, summing over all \( \ell = 1, \ldots, r-1 \) would give
\[ P(\hat{\beta}) = \sum_{\ell=1}^{r-1} \left\| (W^{(\ell)} \ast \hat{\beta})_{g_{r,\ell}} \right\|_2 > \sum_{r=2}^{p} \sum_{\ell=1}^{r-1} \langle W^{(\ell)} \ast \hat{\beta}, \hat{a}^{(\ell)} \rangle, \]
which leads to a contradiction. So \( (\hat{a}^{(\ell)})_{g_{r,\ell}} = \frac{\langle W^{(\ell)} \ast \hat{\beta}, \hat{a}^{(\ell)} \rangle}{\left\| (W^{(\ell)} \ast \hat{\beta})_{g_{r,\ell}} \right\|_2} \) for \( \hat{\beta}_{g_{r,\ell}} \neq 0 \) and \( \left\| \hat{a}^{(\ell)}_{g_{r,\ell}} \right\|_2 \leq 1 \) for \( \hat{\beta}_{g_{r,\ell}} = 0. \quad \Box \)

**K Proof of Lemma 11**

*Proof.* Following from the notation in Appendix F observe that \( \mathcal{L}(\tau, z, \beta; \nu, \phi, a^{(\ell)}) \) is jointly convex in \( \tau, z \) and \( \beta \) and it is strongly convex in \( \tau \) and \( z \). So the minimizer \( \hat{z} \) and \( \hat{\tau} \) are unique.

To see this in more general setting, WLOG suppose \( f(x,y) \) is convex in \( y \) and is strictly convex in \( x \). Then for \( x_1 \neq x_2 \) and \( \theta \in (0, 1) \) we have
\[ f(\theta x_1 + (1-\theta)x_2, y) < \theta f(x_1, y) + (1-\theta)f(x_2, y) \]
Now suppose \((\hat{x}_1, \hat{y})\) and \((\hat{x}_2, \hat{y})\) are both minimum of \( f \), then take \( \theta = 1/2 \) we have \( f\left(\frac{\hat{x}_1 + \hat{x}_2}{2}, \hat{y}\right) < f(\hat{x}_1, \hat{y}) = f(\hat{x}_2, \hat{y}) \), which leads to a contradiction.

By the primal-dual relation, we know that if \( \hat{\beta} \) and \( \bar{\beta} \) are two solutions to (7) in the paper, then \( \hat{\beta}_r = \bar{\beta}_r \) and \( X_{1,r} \hat{\beta} = X_{1,r} \bar{\beta} \). So from the equality \( f(\hat{\beta}) = f(\bar{\beta}) \) we know that \( P(\hat{\beta}) = P(\bar{\beta}) \). Also by
\[ f(\bar{\beta}) = \mathcal{L}(\hat{\tau}, \hat{z}, \bar{\beta}; \hat{\nu}, \hat{\phi}, a^{(\ell)}) \leq \mathcal{L}(\hat{\tau}, \hat{z}, \bar{\beta}; \hat{\nu}, \hat{\phi}, a^{(\ell)}) \leq \mathcal{L}(\hat{\tau}, \hat{z}, \bar{\beta}; \hat{\nu}, \hat{\phi}, a^{(\ell)}) = f(\bar{\beta}), \]

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we have
\[ \mathcal{L}(\hat{\tau}, \hat{\beta}, \hat{\nu}, \hat{\phi}, \hat{\alpha}) = \mathcal{L}(\check{\tau}, \check{\beta}, \check{\nu}, \check{\phi}, \check{\alpha}), \]
and thus
\[ \sum_{\ell=1}^{r-1} \langle W^{(\ell)} \ast \hat{\alpha}, \hat{\beta} \rangle = \sum_{\ell=1}^{r-1} \langle W^{(\ell)} \ast \check{\alpha}, \check{\beta} \rangle = P(\hat{\beta}) = P(\check{\beta}) = \sum_{\ell=1}^{r-1} \left\| W^{(\ell)} \ast \hat{\beta} \right\|_{2}. \]

Now for any \( \ell \leq r - 1 \) suppose \( \left\| (\hat{\alpha})_{g_r, \ell} \right\|_{2} < 1 \), then for the equality above to hold, we must have \( \hat{\beta}_{g_r, \ell} = 0 \). Therefore, by Lemma 10, \( \hat{\beta}_{g_r, \ell} = 0 \implies \hat{\beta}_{g_r, \ell} = 0 \), so any other solutions to (7) in the paper cannot be less sparse than \( \hat{\beta} \).

\[ \]
Next we deal with the second term in $M_n$, recall from (33) that
\[
\frac{4}{n^2 \lambda^2} \mathbb{E}_R [O_T E_r]_2^2 = \frac{4}{n^2} \left( \frac{1}{2} X_r^T C_I + \sqrt{\frac{1}{4} (X_r^T C_s)^2 + \frac{4}{n} \mathbb{E}_R [O_T E_r]_2^2} \right)^2 \mathbb{E}_R [O_T E_r]_2^2 \\
\leq \frac{4}{n^2} \left( \frac{1}{4} (X_r^T C_s)^2 + \frac{4}{n} \mathbb{E}_R [O_T E_r]_2^2 \right) \mathbb{E}_R [O_T E_r]_2^2 \\
= \frac{(X_r^T C_s)^2}{\mathbb{E}_R [O_T E_r]_2^2} + \frac{16}{\lambda^2}.
\]

The next lemma gives us a handle on the numerator of the first term.

**Lemma 19.** Using the general weight (6) in the paper, conditioned on $X_I$, we have
\[
P \left[ |X_r^T C_s| \geq 1 \right] \leq 2 \exp \left( -\frac{n \alpha^2}{3 \theta^2 r^2 \pi^2 K} \right) + 2 \exp \left( -\frac{n}{2} \right).
\]

**Proof.** Conditioned on $X_I$, from the decomposition (35) and the definition of $C_I$
\[
X_r^T C_I = \Sigma_r \Sigma_{rI}^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right) + E_r^T X_I (X_r^T X_I)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right)_I.
\]
By irrepresentable condition [A3 and 37]
\[
\Sigma_r \Sigma_{rI}^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right)_I \leq 1 - \alpha.
\]
Note that since $\text{Var} (E_{ir}) = \theta^2_i$ for $i = 1, \ldots, n$. Let $B^{(r)} := E_r^T X_I (X_r^T X_I)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right)_I$, then by Lemma 18 $B^{(r)}$ has mean zero and variance at most
\[
\text{Var} \left( B^{(r)} \mid X_I \right) = \frac{\theta^2_i}{n} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right)_I^T \left( \frac{1}{n} X_r^T X_I \right)^{-1} \left( \sum_{\ell=1}^{r-1} W^{(\ell)} * \tilde{a}^{(\ell)} \right)_I \leq \frac{3 \theta^2_i \kappa^2 \pi^2 K}{2n},
\]
with probability greater than $1 - 2 \exp \left( \frac{n}{2} \right)$. Then by Lemma 13, we have the
\[
P \left[ B^{(r)} \geq \alpha \right] \leq 2 \exp \left( -\frac{n \alpha^2}{3 \theta^2_i \kappa^2 \pi^2 K} \right) + 2 \exp \left( -\frac{n}{2} \right).
\]

Since $\|O_T E_r\|_2^2 \sim \chi^2 (n - K)$. To bound it, we cite a concentration inequality from Wainwright (2009) (in specific (54b)) as the following lemma:

**Lemma 20** (Tail Bounds for $\chi^2$-variates). For a centralized $\chi^2$-variate $X$ with $d$-degrees of freedom, for all $\varepsilon \in (0, 1/2)$, we have
\[
P [X \leq d(1 - \varepsilon)] \leq \exp \left( -\frac{1}{4} d \varepsilon^2 \right).
\]
From lemma 20 it follows that

$$P \left[ \|O_T E_r\|^2 \leq \theta_r \left( n - K \right) (1 - \varepsilon) \right] \leq \exp \left( -\frac{1}{4} \left( n - K \right) \varepsilon^2 \right),$$

which together with lemma 19 imply that

$$P \left[ \left( X^T C_T \right)^2 \geq \frac{1}{\theta_r \left( n - K \right) (1 - \varepsilon)} \right] \leq 2 \exp \left( -\frac{na^2}{3\theta_r \kappa^2 n^2 K} \right) + 2 \exp \left( -\frac{n}{2} \right) + \exp \left( -\frac{1}{4} \left( n - K \right) \varepsilon^2 \right).$$

The result follows from a union bound.

## N Proof of Lemma 15

**Proof.** For the design matrix $X_{n \times p}$ with independent rows, denote $X_i = (X_i)^T \in \mathbb{R}^p$. Then $X_i$ are i.i.d with mean 0 and true covariance matrix $\Sigma = (L^T L)^{-1}$ for $i = 1, ..., n$. And $\bar{X} = \frac{1}{n} \sum_{i=1}^n X_i$ has mean 0 and true covariance matrix $\frac{1}{n} \Sigma$.

Let $Y_i = L X_i \in \mathbb{R}^p$. Then $Y_i$ are i.i.d with mean 0 and true covariance matrix $L \Sigma L^T = L \left( L^T L \right)^{-1} L^T = I_p$. And $\bar{Y} = \frac{1}{n} \sum_{i=1}^n Y_i = \frac{1}{n} \sum_{i=1}^n L X_i = L \bar{X}$ has mean zero and covariance matrix $\frac{1}{n} I_p$. Also the corresponding design matrix $Y = XL^T$ has independent rows.

$$SL^T = \frac{1}{n} \sum_{i=1}^n X_i (X_i - \bar{X})^T, \quad L^T = \frac{1}{n} \sum_{i=1}^n (X_i - \bar{X}) (L X_i - L \bar{X})^T = \frac{1}{n} \sum_{i=1}^n (X_i - \bar{X}) (Y_i - \bar{Y})^T.$$

So we have

$$(SL^T)_{ij} = n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - \bar{X}_i \bar{Y}_j.$$

Recall that

$$W = SL^T - (L)^{-1},$$

then

$$|W_{ij}| \leq \left| n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - (L^T)^{-1}_{ij} \right| + |\bar{X}_i \bar{Y}_j|.$$

$$P \left[ \max_{ij} |W_{ij}| > t \right] \leq P \left[ \max_{ij} \left| n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - (L)^{-1}_{ij} \right| > \frac{t}{2} \right] + P \left[ \max_{ij} |\bar{X}_i \bar{Y}_j| > \frac{t}{2} \right]$$

$$\leq P \left[ \left| n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - (L)^{-1}_{ij} \right| > \frac{t}{2} \right] + \sum_{ij} P \left[ \left| n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - (L)^{-1}_{ij} \right| > \frac{t}{2} \right] + \sum_{i} P \left[ |\bar{X}_i| > \sqrt{\frac{t}{2}} \right] + \sum_{j} P \left[ |\bar{Y}_j| > \sqrt{\frac{t}{2}} \right]$$

$$\leq p^2 \max_{ij} P \left[ \left| n^{-1} \sum_{k=1}^p X_{ki} Y_{kj} - (L)^{-1}_{ij} \right| > \frac{t}{2} \right] + p \max_{i} P \left[ |\bar{X}_i| > \sqrt{\frac{t}{2}} \right] + p \max_{j} P \left[ |\bar{Y}_j| > \sqrt{\frac{t}{2}} \right]$$

$$:= p^2 \max_{ij} I_{ij} + p \max_{i} I_{i}^X + p \max_{j} I_{j}^Y.$$

Consider $I_{i}^X$ first. Since $X_{ki}$ are independent sub-Gaussian with variance $(\Sigma)_{ii}$ for $k = 1, .., n$, we have

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\[
E \exp \left( t \frac{\bar{X}_i}{\sqrt{\Sigma_{ii}/n}} \right) = \prod_{k=1}^{n} E \exp \left( t \frac{X_{ki}}{\sqrt{n \Sigma_{ii}}} \right) \quad \text{by independence}
\]
\[
\leq \prod_{k=1}^{n} \exp \left( \tilde{C}_1 t^2/n \right) = \exp(\tilde{C}_1 t^2) \quad \text{by the definition of sub-Gaussian},
\]
so \( \bar{X}_i \) is sub-Gaussian with variance \( \Sigma_{ii}/n \).

By lemma 5.5 in [Vershynin (2010)], we have
\[
P \left[ \left| \bar{X}_i \right| / \sqrt{\Sigma_{ii}} > t \right] \leq \exp \left( 1 - \frac{t^2}{K^2} \right),
\]
where \( K_1 \) is a constant that does not depend on \( i \).

Following the same argument we have
\[
E \exp \left( t \frac{\bar{Y}_i}{\sqrt{1/n}} \right) = \prod_{k=1}^{n} E \exp \left( t \frac{Y_{ki}}{\sqrt{n}} \right) \leq \exp \left( \tilde{C}_2 t^2 \right),
\]
thus
\[
P \left[ \left| \bar{Y}_i \right| / \sqrt{1/n} > t \right] \leq \exp \left( 1 - \frac{t^2}{K_2^2} \right),
\]
where \( K_2 \) is a constant that does not depend on \( i \). And we have
\[
I_i^X + I_i^Y = P \left[ \left| \bar{X}_i \right| > \sqrt{t/2} \right] + P \left[ \left| \bar{Y}_i \right| > \sqrt{t/2} \right]
\]
\[
= P \left[ \frac{\left| \bar{X}_i \right|}{\sqrt{\Sigma_{ii}/n}} > \sqrt{t/2} \right] + P \left[ \frac{\left| \bar{Y}_i \right|}{\sqrt{1/n}} > \sqrt{t/2} \right] \leq \exp \left( 1 - \frac{nt}{2K^2 \Sigma_{ii}} \right) + \exp \left( 1 - \frac{nt}{2K_2^2} \right).
\]

Thus
\[
\max_{i} (I_i^X + I_i^Y) \leq 8 \exp \left( - \frac{C_1 nt}{\max_{i} \Sigma_{ii}^*} \right),
\]
for some constant \( C_1 \).

Now consider the term \( I_{ij} \). We have shown that both \( X \) and \( Y \) have independent rows. So for any \( i, j, Z_{k}^{(ij)} = X_{ki}Y_{kj} \) are independent for \( k = 1, \ldots, n \). Note that
\[
E (X_{ki}Y_{kj}) = \text{Cov} (X, LX)_{ij} = 0 = \left[ \text{Cov} (X, X) L^T \right]_{ij} = (\Sigma L^T)_{ij} = (L)^{-1}_{ij}.
\]

If there exist \( \nu_{ij} \) and \( c_{ij} \) such that
\[
\sum_{k=1}^{n} E \left( X_{ki}^2 Y_{kj}^2 \right) \leq \nu_{ij}
\]
\[
\sum_{k=1}^{n} E \left\{ \left( X_{ki}Y_{kj} \right)^q \right\} = \frac{q!}{2} \nu_{ij} c_{ij}^{q-2} \quad \text{for some} \quad q \geq 3 \in \mathbb{N},
\]
then by Theorem 2.10 (Corollary 2.11) in [Boucheron et al. (2013)], \( \forall t > 0 \), we have
\[
P \left[ \sum_{k=1}^{n} \left( X_{ki}Y_{kj} - (L)^{-1}_{ij} \right) > t \right] \leq 2 \exp \left( - \frac{t^2}{2(\nu_{ij} + c_{ij} t)} \right).
\]

To find \( \nu_{ij} \), by Lemma 5.5 in [Vershynin (2010)]
\[
\left( E \left| X_{ij} / \sqrt{\Sigma_{jj}} \right|^q \right)^{1/q} \leq K_3 \sqrt{q},
\]

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for all \( q \geq 1 \) and some constant \( K_3 \) that does not depend on \( j \), which implies
\[
E |X_{ij}|^q \leq K_3^q q^{q/2} (\Sigma_{ii})^{q/2}.
\]

Similarly,
\[
E |Y_{ij}|^q \leq K_4^q q^{q/2},
\]
for all \( q \geq 1 \) and some constant \( K_4 \) that does not depend on \( j \). Therefore,
\[
\sum_{k=1}^{n} E \left( X_{ki}^2 Y_{kj}^2 \right) \leq \sum_{k=1}^{n} \sqrt{E X_{ki}^4 E Y_{kj}^4} \leq n \sqrt{K_3^4 K_4^4 (\Sigma_{ii})^2} = 16n K_3^2 K_4^2 \Sigma_{ii}.
\]

Similarly,
\[
\sum_{k=1}^{n} E \left\{ (X_{ki} Y_{kj})^q \right\} \leq \sum_{k=1}^{n} \sqrt{E X_{ki}^{2q} E Y_{kj}^{2q}} \leq n \sqrt{K_3^{2q} (2q)^2 K_4^{2q} (\Sigma_{ii})^2} = n K_3^q K_4^q (2q)^q (\Sigma_{ii})^{q/2}.
\]
So taking
\[
\nu_{ij} = K_5 n \Sigma_{ii}^* \\
c_{ij} = K_5 \sqrt{\Sigma_{ii}^*}
\]
for some \( K_5 \) large enough and does not depend on \( i,j \).

Now we have
\[
I_{ij} \leq 2 \exp \left( -\frac{n^2 t^2}{4 (2 \nu_{ij} + c_{ij} nt)} \right) = 2 \exp \left( -\frac{nt^2}{4 (2 K_5 \Sigma_{ii}^* + K_5 \sqrt{\Sigma_{ii}} t)} \right).
\]
If \( t \leq 2 \max_i \sqrt{\Sigma_{ii}^*} \), then with \( C_2 = (16K_5)^{-1} \) we have
\[
I_{ij} \leq 2 \exp \left( -\frac{C_2 n t^2}{\max_i \Sigma_{ii}^*} \right).
\]

To sum up, for any \( 0 < t \leq 2 \max_i \sqrt{\Sigma_{ii}^*} \),
\[
P \left[ \max_{ij} |W_{ij}| > t \right] \leq 2 p^2 \exp \left( -\frac{C_2 n t^2}{\max_i \Sigma_{ii}^*} \right) + 8p \exp \left( -\frac{C_1 n t}{\max_i \Sigma_{ii}^*} \right).
\]

\[\square\]

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