

Confidence Intervals and Hypothesis Testing for High-Dimensional Regression

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April 18, 2014

Abstract

Fitting high-dimensional statistical models often requires the use of non-linear parameter estimation procedures. As a consequence, it is generally impossible to obtain an exact characterization of the probability distribution of the parameter estimates. This in turn implies that it is extremely challenging to quantify the *uncertainty* associated with a certain parameter estimate. Concretely, no commonly accepted procedure exists for computing classical measures of uncertainty and statistical significance as confidence intervals or p -values for these models.

We consider here high-dimensional linear regression problem, and propose an efficient algorithm for constructing confidence intervals and p -values. The resulting confidence intervals have nearly optimal size. When testing for the null hypothesis that a certain parameter is vanishing, our method has nearly optimal power.

Our approach is based on constructing a ‘de-biased’ version of regularized M-estimators. The new construction improves over recent work in the field in that it does not assume a special structure on the design matrix. We test our method on synthetic data and a high-throughput genomic data set about riboflavin production rate, made publicly available by [BKM14].

1 Introduction

It is widely recognized that modern statistical problems are increasingly high-dimensional, i.e. require estimation of more parameters than the number of observations/samples. Examples abound from signal processing [LDSP08], to genomics [PZB⁺10], collaborative filtering [KBV09] and so on. A number of successful estimation techniques have been developed over the last ten years to tackle these problems. A widely applicable approach consists in optimizing a suitably regularized likelihood function. Such estimators are, by necessity, non-linear and non-explicit (they are solution of certain optimization problems).

The use of non-linear parameter estimators comes at a price. In general, it is impossible to characterize the distribution of the estimator. This situation is very different from the one of classical statistics in which either exact characterizations are available, or asymptotically exact ones can be derived from large sample theory [VdV00]. This has an important and very concrete consequence.

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In classical statistics, generic and well accepted procedures are available for characterizing the uncertainty associated to a certain parameter estimate in terms of confidence intervals or p -values [Was04, LR05]. However, no analogous procedures exist in high-dimensional statistics.

In this paper we develop a computationally efficient procedure for constructing confidence intervals and p -values for a broad class of high-dimensional regression problems. The salient features of our procedure are:

- (i) Our approach guarantees nearly optimal confidence interval sizes and testing power.
- (ii) It is the first one to achieve this goal under essentially no assumptions beyond the standard conditions for high-dimensional consistency.
- (iii) It allows for a streamlined analysis with respect to earlier work in the same area.

For the sake of clarity, we will focus our presentation on the case of linear regression, under Gaussian noise. Section 4 provides a detailed study of the case of non-Gaussian noise. A preliminary report on our results was presented in NIPS 2013 [JM13a], which also discusses generalizations of the same approach to generalized linear models, and regularized maximum likelihood estimation.

In a linear regression model, we are given n i.i.d. pairs $(Y_1, X_1), (Y_2, X_2), \dots, (Y_n, X_n)$, with vectors $X_i \in \mathbb{R}^p$ and response variables Y_i given by

$$Y_i = \langle \theta_0, X_i \rangle + W_i, \quad W_i \sim \mathbf{N}(0, \sigma^2). \quad (1)$$

Here $\theta_0 \in \mathbb{R}^p$ and $\langle \cdot, \cdot \rangle$ is the standard scalar product in \mathbb{R}^p . In matrix form, letting $Y = (Y_1, \dots, Y_n)^\top$ and denoting by \mathbf{X} the design matrix with rows $X_1^\top, \dots, X_n^\top$, we have

$$Y = \mathbf{X} \theta_0 + W, \quad W \sim \mathbf{N}(0, \sigma^2 \mathbf{I}_{n \times n}). \quad (2)$$

The goal is to estimate the unknown (but fixed) vector of parameters $\theta_0 \in \mathbb{R}^p$.

In the classic setting, $n \gg p$ and the estimation method of choice is ordinary least squares yielding $\hat{\theta}^{\text{OLS}} = (\mathbf{X}^\top \mathbf{X})^{-1} \mathbf{X}^\top Y$. In particular $\hat{\theta}^{\text{OLS}}$ is Gaussian with mean θ_0 and covariance $\sigma^2 (\mathbf{X}^\top \mathbf{X})^{-1}$. This directly allows to construct confidence intervals¹.

In the high-dimensional setting where $p > n$, the matrix $(\mathbf{X}^\top \mathbf{X})$ is rank deficient and one has to resort to biased estimators. A particularly successful approach is the LASSO [Tib96, CD95] which promotes sparse reconstructions through an ℓ_1 penalty:

$$\hat{\theta}^n(Y, \mathbf{X}; \lambda) \equiv \arg \min_{\theta \in \mathbb{R}^p} \left\{ \frac{1}{2n} \|Y - \mathbf{X}\theta\|_2^2 + \lambda \|\theta\|_1 \right\}. \quad (3)$$

In case the right hand side has more than one minimizer, one of them can be selected arbitrarily for our purposes. We will often omit the arguments Y, \mathbf{X} , as they are clear from the context.

We denote by $S \equiv \text{supp}(\theta_0)$ the support of $\theta_0 \in \mathbb{R}^p$, defined as

$$\text{supp}(\theta_0) \equiv \{i \in [p] : \theta_{0,i} \neq 0\},$$

where we use the notation $[p] = \{1, \dots, p\}$. We further let $s_0 \equiv |S|$. A copious theoretical literature [CT05, BRT09, BvdG11] shows that, under suitable assumptions on \mathbf{X} , the LASSO is nearly as

Table 1: Unbiased estimator for θ_0 in high-dimensional linear regression models

Input: Measurement vector y , design matrix \mathbf{X} , parameters λ, μ .

Output: Unbiased estimator $\hat{\theta}^u$.

- 1: Let $\hat{\theta}^n = \hat{\theta}^n(Y, \mathbf{X}; \lambda)$ be the LASSO estimator as per Eq. (3).
- 2: Set $\hat{\Sigma} \equiv (\mathbf{X}^\top \mathbf{X})/n$.
- 3: **for** $i = 1, 2, \dots, p$ **do**
- 4: Let m_i be a solution of the convex program:

$$\begin{aligned} & \text{minimize} && m^\top \hat{\Sigma} m \\ & \text{subject to} && \|\hat{\Sigma} m - e_i\|_\infty \leq \mu, \end{aligned} \tag{4}$$

where $e_i \in \mathbb{R}^p$ is the vector with one at the i -th position and zero everywhere else.

- 5: Set $M = (m_1, \dots, m_p)^\top$. If any of the above problems is not feasible, then set $M = \mathbf{I}_{p \times p}$.
- 6: Define the estimator $\hat{\theta}^u$ as follows:

$$\hat{\theta}^u = \hat{\theta}^n(\lambda) + \frac{1}{n} M \mathbf{X}^\top (Y - \mathbf{X} \hat{\theta}^n(\lambda)) \tag{5}$$

accurate as if the support S was known *a priori*. Namely, for $n = \Omega(s_0 \log p)$, we have $\|\hat{\theta}^n - \theta_0\|_2^2 = O(s_0 \sigma^2 (\log p)/n)$.

As mentioned above, these remarkable properties come at a price. Deriving an exact characterization for the distribution of $\hat{\theta}^n$ is not tractable in general, and hence there is no simple procedure to construct confidence intervals and p -values. A closely related property is that $\hat{\theta}^n$ is biased, an unavoidable property in high dimension, since a point estimate $\hat{\theta}^n \in \mathbb{R}^p$ must be produced from data in lower dimension $Y \in \mathbb{R}^n$, $n < p$. We refer to Section 2.2 for further discussion of this point.

In order to overcome this challenge, we construct a de-biased estimator from the LASSO solution. The de-biased estimator is given by the simple formula $\hat{\theta}^u = \hat{\theta}^n + (1/n) M \mathbf{X}^\top (Y - \mathbf{X} \hat{\theta}^n)$, as in Eq. (5). The basic intuition is that $\mathbf{X}^\top (Y - \mathbf{X} \hat{\theta}^n)/(n\lambda)$ is a subgradient of the ℓ_1 norm at the LASSO solution $\hat{\theta}^n$. By adding a term proportional to this subgradient, our procedure compensates the bias introduced by the ℓ_1 penalty in the LASSO.

We will prove in Section 2.1 that $\hat{\theta}^u$ is approximately Gaussian, with mean θ_0 and covariance $\sigma^2(M \hat{\Sigma} M)/n$, where $\hat{\Sigma} = (\mathbf{X}^\top \mathbf{X}/n)$ is the empirical covariance of the feature vectors. This result allows to construct confidence intervals and p -values in complete analogy with classical statistics procedures. For instance, letting $Q \equiv M \hat{\Sigma} M$, $[\hat{\theta}_i^u - 1.96\sigma \sqrt{Q_{ii}/n}, \hat{\theta}_i^u + 1.96\sigma \sqrt{Q_{ii}/n}]$ is a 95% confidence interval. The size of this interval is of order σ/\sqrt{n} , which is the optimal (minimum) one, i.e. the same that would have been obtained by knowing *a priori* the support of θ_0 . In practice the noise standard deviation is not known, but σ can be replaced by any consistent estimator $\hat{\sigma}$ (see Section 3 for more details on this).

A key role is played by the matrix $M \in \mathbb{R}^{p \times p}$ whose function is to ‘decorrelate’ the columns of \mathbf{X} . We propose here to construct M by solving a convex program that aims at optimizing two

¹For instance, letting $Q \equiv (\mathbf{X}^\top \mathbf{X}/n)^{-1}$, $[\hat{\theta}_i^{\text{OLS}} - 1.96\sigma \sqrt{Q_{ii}/n}, \hat{\theta}_i^{\text{OLS}} + 1.96\sigma \sqrt{Q_{ii}/n}]$ is a 95% confidence interval [Was04].

objectives. On one hand, we try to control $|M\widehat{\Sigma} - I|_\infty$ (here and below $|\cdot|_\infty$ denotes the entrywise ℓ_∞ norm) which –as shown in Theorem 2.5– controls the non-Gaussianity and bias of $\widehat{\theta}^u$. On the other, we minimize $[M\widehat{\Sigma}M]_{i,i}$, for each $i \in [p]$, which controls the variance of $\widehat{\theta}_i^u$.

The idea of constructing a de-biased estimator of the form $\widehat{\theta}^u = \widehat{\theta}^n + (1/n)M\mathbf{X}^\top(Y - \mathbf{X}\widehat{\theta}^n)$ was used by the present authors in [JM13b], that suggested the choice $M = c\Sigma^{-1}$, with $\Sigma = \mathbb{E}\{X_1X_1^\top\}$ the population covariance matrix and c a positive constant. A simple estimator for Σ was proposed for sparse covariances, but asymptotic validity and optimality were proven only for uncorrelated Gaussian designs (i.e. Gaussian \mathbf{X} with $\Sigma = I$). Van de Geer, Bühlmann, Ritov and Dezeure [vdGBRD13] used the same construction with M an estimate of Σ^{-1} which is appropriate for sparse inverse covariances. These authors prove semi-parametric optimality in a non-asymptotic setting, provided the sample size is at least $n = \Omega((s_0 \log p)^2)$.

From a technical point of view, our proof starts from a simple decomposition of the de-biased estimator $\widehat{\theta}^u$ into a Gaussian part and an error term, already used in [vdGBRD13]. However –departing radically from earlier work– we realize that M need not be a good estimator of Σ^{-1} in order for the de-biasing procedure to work. We instead set M as to minimize the error term and the variance of the Gaussian term. As a consequence of this choice, our approach applies to general covariance structures Σ . By contrast, earlier approaches applied only to sparse Σ , as in [JM13b], or sparse Σ^{-1} as in [vdGBRD13]. The only assumptions we make on Σ are the standard compatibility conditions required for high-dimensional consistency [BvdG11]. A detailed comparison of our results with the ones of [vdGBRD13] can be found in Section 2.3.

Our presentation is organized as follows.

Section 2 considers a general debiased estimator of the form $\widehat{\theta}^u = \widehat{\theta}^n + (1/n)M\mathbf{X}^\top(Y - \mathbf{X}\widehat{\theta}^n)$.

We introduce a figure of merit of the pair M, \mathbf{X} , termed the generalized coherence parameter $\mu_*(\mathbf{X}; M)$. We show that, if the generalized coherence is small, then the debiasing procedure is effective (for a given deterministic design), see Theorem 2.3.

We then turn to random designs, and show that the generalized coherence parameter can be made as small as $\sqrt{(\log p)/n}$, though a convex optimization procedure for computing M . This results in a bound on the bias of $\widehat{\theta}^u$, cf. Theorem 2.5: the largest entry of the bias is of order $(s_0 \log p)/n$. This must be compared with the standard deviation of $\widehat{\theta}_i^u$, which is of order σ/\sqrt{n} . The conclusion is that, for $s_0 = o(\sqrt{n}/\log p)$, the bias of $\widehat{\theta}^u$ is negligible.

Section 3 applies these distributional results to deriving confidence intervals and hypothesis testing procedures for low-dimensional marginals of $\widehat{\theta}_0$. The basic intuition is that $\widehat{\theta}^u$ is approximately Gaussian with mean θ_0 , and known covariance structure. Hence standard optimal tests can be applied.

We prove a general lower bound on the power of our testing procedure, in Theorem 3.5. In the special case of Gaussian random designs with i.i.d. rows, we can compare this with the upper bound proved in [JM13b], cf. Theorem 3.6. As a consequence, the asymptotic efficiency of our approach is constant-optimal. Namely, it is lower bounded by a constant $1/\eta_{\Sigma, s_0}$ which is bounded away from 0, cf. Theorem 3.7. (For instance $\eta_{I, s_0} = 1$, and η_{Σ, s_0} is always upper bounded by the condition number of Σ .)

Section 4 uses the central limit theorem for triangular arrays to generalize the above results to non-Gaussian noise.

Section 5 illustrates the above results through numerical simulations both on synthetic and on real data. In the interest of reproducibility, an R implementation of our algorithm is available at <http://www.stanford.edu/~montanar/sslasso/>.

Note that our proofs require stricter sparsity s_0 (or larger sample size n) than required for consistent estimation. We assume $s_0 = o(\sqrt{n}/\log p)$ instead of $s_0 = o(n/\log p)$ [CT07, BRT09, BvdG11]. The same assumption is made in [vdGBRD13], on top of additional assumptions on the sparsity of Σ^{-1} .

It is currently an open question whether successful hypothesis testing can be performed under the weaker assumption $s_0 = o(n/\log p)$. We refer to [JM13c] for preliminary work in that direction. The barrier at $s_0 = o(\sqrt{n}/\log p)$ is possibly related to an analogous assumption that arises in Gaussian graphical models selection [RSZZ13].

1.1 Further related work

The theoretical literature on high-dimensional statistical models is vast and rapidly growing. Estimating sparse linear regression models is the most studied problem in this area, and a source of many fruitful ideas. Limiting ourselves to linear regression, earlier work investigated prediction error [GR04], model selection properties [MB06, ZY06, Wai09, CP09], ℓ_2 consistency [CT05, BRT09]. . Of necessity, we do not provide a complete set of references, and instead refer the reader to [BvdG11] for an in-depth introduction to this area.

The problem of quantifying statistical significance in high-dimensional parameter estimation is, by comparison, far less understood. Zhang and Zhang [ZZ14], and Bühlmann [Büh13] proposed hypothesis testing procedures under restricted eigenvalue or compatibility conditions [BvdG11]. These papers provide deterministic guarantees but –in order to achieve a certain target significance level α and power $1 - \beta$ – they require $|\theta_{0,i}| \geq c \max\{\sigma s_0 \log p/n, \sigma/\sqrt{n}\}$. The best lower bound [JM13b] shows that any such test requires instead $|\theta_{0,i}| \geq c(\alpha, \beta)\sigma/\sqrt{n}$. (The lower bound of [JM13b] is reproduced as Theorem 3.6 here, for the reader’s convenience.)

In other words, the guarantees of [ZZ14, Büh13] can be suboptimal by a factor as large as $\sqrt{s_0}$. Equivalently, in order for the coefficient $\theta_{0,i}$ to be detectable with appreciable probability, it needs to be larger than the overall ℓ_2 error. Here we will propose a test that –for random designs– achieves significance level α and power $1 - \beta$ for $|\theta_{0,i}| \geq c'(\alpha, \beta)\sigma/\sqrt{n}$.

Lockhart et al. [LTTT13] develop a test for the hypothesis that a newly added coefficient along the LASSO regularization path is irrelevant. This however does not allow to test arbitrary coefficients at a given value of λ , which is instead the problem addressed in this paper. These authors further assume that the current LASSO support contains the actual support $\text{supp}(\theta_0)$ and that the latter has bounded size.

Belloni, Chernozhukov and collaborators [BCH11, BCW13] consider inference in a regression model with high-dimensional data. In this model the response variable relates to a scalar main regressor and a p -dimensional control vector. The main regressor is of primary interest and the control vector is treated as nuisance component. Assuming that the control vector is s_0 -sparse, the authors propose a method to construct confidence regions for the parameter of interest under the sample size requirement $(s_0^2 \log p)/n \rightarrow 0$. The proposed method is shown to attain the semi-parametric efficiency bounds for this class of models. The key modeling assumption in this paper is that the scalar regressor of interest is random, and depends linearly on the p -dimensional control vector, with a sparse coefficient vector (with sparsity again of order $o(\sqrt{n/\log p})$). This assumption

is closely related to the sparse inverse covariance assumption of [vdGBRD13] (with the difference that only one regressor is tested).

Finally, resampling methods for hypothesis testing were studied in [MB10, MTC11]. These methods are perturbation-based procedures to approximate the distribution of a general class of penalized parameter estimates for the case $n > p$. The idea is to consider the minimizer of a stochastically perturbed version of the regularized objective function, call it $\tilde{\theta}$, and characterize the limiting distribution of the regularized estimator $\hat{\theta}$ in terms of the distribution of $\tilde{\theta}$. In order to estimate the latter, a large number of random samples of the perturbed objective function are generated, and for each sample the minimizer is computed. Finally the theoretical distribution of $\tilde{\theta}$ is approximated by the empirical distribution of these minimizers.

After the present paper was submitted for publication, we became aware that Bühlmann and Dezeure [DB13] had independently worked on similar ideas.

1.2 Preliminaries and notations

In this section we introduce some basic definitions used throughout the paper, starting with simple notations.

For a matrix A and set of indices I, J , we let $A_{I,J}$ denote the submatrix formed by the rows in I and columns in J . Also, $A_{I,\cdot}$ (resp. $A_{\cdot,J}$) denotes the submatrix containing just the rows (reps. columns) in I . Likewise, for a vector v , v_I is the restriction of v to indices in I . We use the shorthand $A_{I,J}^{-1} = (A^{-1})_{I,J}$. In particular, $A_{i,i}^{-1} = (A^{-1})_{i,i}$. The maximum and the minimum singular values of A are respectively denoted by $\sigma_{\max}(A)$ and $\sigma_{\min}(A)$. We write $\|v\|_p$ for the standard ℓ_p norm of a vector v , i.e., $\|v\|_p = (\sum_i |v_i|^p)^{1/p}$. and $\|v\|_0$ for the number of nonzero entries of v . For a matrix A , $\|A\|_p$ is the ℓ_p operator norm, and $|A|_p$ is the elementwise ℓ_p norm. For a vector v , $\text{supp}(v)$ represents the positions of nonzero entries of v . Throughout, $\Phi(x) \equiv \int_{-\infty}^x e^{-t^2/2} dt / \sqrt{2\pi}$ denotes the CDF of the standard normal distribution. Finally, *with high probability* (w.h.p) means with probability converging to one as $n \rightarrow \infty$.

We let $\hat{\Sigma} \equiv \mathbf{X}^\top \mathbf{X} / n$ be the sample covariance matrix. For $p > n$, $\hat{\Sigma}$ is always singular. However, we may require $\hat{\Sigma}$ to be nonsingular for a restricted set of directions.

Definition 1.1. *Given a symmetric matrix $\hat{\Sigma} \in \mathbb{R}^{p \times p}$ and a set $S \subseteq [p]$, the corresponding compatibility constant is defined as*

$$\phi^2(\hat{\Sigma}, S) \equiv \min_{\theta \in \mathbb{R}^p} \left\{ \frac{|S| \langle \theta, \hat{\Sigma} \theta \rangle}{\|\theta_S\|_1^2} : \theta \in \mathbb{R}^p, \|\theta_{S^c}\|_1 \leq 3\|\theta_S\|_1 \right\}. \quad (6)$$

We say that $\hat{\Sigma} \in \mathbb{R}^{p \times p}$ satisfies the compatibility condition for the set $S \subseteq [p]$, with constant ϕ_0 if $\phi(\hat{\Sigma}, S) \geq \phi_0$. We say that it holds for the design matrix \mathbf{X} , if it holds for $\hat{\Sigma} = \mathbf{X}^\top \mathbf{X} / n$.

In the following, we shall drop the argument $\hat{\Sigma}$ if clear from the context. Note that a slightly more general definition is used normally [BvdG11, Section 6.13], whereby the condition $\|\theta_{S^c}\|_1 \leq 3\|\theta_S\|_1$, is replaced by $\|\theta_{S^c}\|_1 \leq L\|\theta_S\|_1$. The resulting constant $\phi(\hat{\Sigma}, S, L)$ depends on L . For the sake of simplicity, we restrict ourselves to the case $L = 3$.

Definition 1.2. *The sub-gaussian norm of a random variable X , denoted by $\|X\|_{\psi_2}$, is defined as*

$$\|X\|_{\psi_2} = \sup_{q \geq 1} q^{-1/2} (\mathbb{E}|X|^q)^{1/q}.$$

For a random vector $X \in \mathbb{R}^n$, its sub-gaussian norm is defined as $\|X\|_{\psi_2} = \sup_{x \in S^{n-1}} \|\langle X, x \rangle\|_{\psi_2}$, where S^{n-1} denotes the unit sphere in \mathbb{R}^n .

Definition 1.3. The sub-exponential norm of a random variable X , denoted by $\|X\|_{\psi_1}$, is defined as

$$\|X\|_{\psi_1} = \sup_{q \geq 1} q^{-1} (\mathbb{E}|X|^q)^{1/q}.$$

For a random vector $X \in \mathbb{R}^n$, its sub-exponential norm is defined as $\|X\|_{\psi_1} = \sup_{x \in S^{n-1}} \|\langle X, x \rangle\|_{\psi_1}$, where S^{n-1} denotes the unit sphere in \mathbb{R}^n .

2 Compensating the bias of the LASSO

In this section we present our characterization of the de-biased estimator $\hat{\theta}^u$ (subsection 2.1). This characterization also clarifies in what sense the LASSO estimator is biased. We discuss this point in subsection 2.2.

2.1 A de-biased estimator for θ_0

As emphasized above, our approach is based on a de-biased estimator defined in Eq. (5), and on its distributional properties. In order to clarify the latter, it is convenient to begin with a slightly broader setting and consider a general debiasing procedure that makes use of an arbitrary $M \in \mathbb{R}^{p \times p}$. Namely, we define

$$\hat{\theta}^*(Y, \mathbf{X}; M, \lambda) = \hat{\theta}^n(\lambda) + \frac{1}{n} M \mathbf{X}^\top (Y - \mathbf{X} \hat{\theta}^n(\lambda)). \quad (7)$$

For notational simplicity, we shall omit the arguments $Y, \mathbf{X}, M, \lambda$ unless they are required for clarity. The quality of this debiasing procedure depends of course on the choice of M , as well as on the design \mathbf{X} . We characterize the pair (\mathbf{X}, M) by the following figure of merit.

Definition 2.1. Given the pair $\mathbf{X} \in \mathbb{R}^{n \times p}$ and $M \in \mathbb{R}^{p \times p}$, let $\hat{\Sigma} = \mathbf{X}^\top \mathbf{X} / n$ denote the associated sample covariance. Then, the generalized coherence parameter of \mathbf{X}, M , denoted by $\mu_*(\mathbf{X}; M)$, is

$$\mu_*(\mathbf{X}; M) \equiv |M \hat{\Sigma} - \mathbf{I}|_\infty. \quad (8)$$

The minimum (generalized) coherence of \mathbf{X} is $\mu_{\min}(\mathbf{X}) = \min_{M \in \mathbb{R}^{p \times p}} \mu_*(\mathbf{X}; M)$. We denote by $M_{\min}(\mathbf{X})$ any minimizer of $\mu_*(\mathbf{X}; M)$.

Note that the minimum coherence can be computed efficiently since $M \mapsto \mu_*(\mathbf{X}; M)$ is a convex function (even more, the optimization problem is a linear program).

The motivation for our terminology can be grasped by considering the following special case.

Remark 2.2. Assume that the columns of \mathbf{X} are normalized to have ℓ_2 norm equal to \sqrt{n} (i.e. $\|\mathbf{X}e_i\|_2 = \sqrt{n}$ for all $i \in [p]$), and $M = \mathbf{I}$. Then $(M \hat{\Sigma} - \mathbf{I})_{i,i} = 0$, and the maximum $|M \hat{\Sigma} - \mathbf{I}|_\infty = \max_{i \neq j} |(\hat{\Sigma})_{ij}|$. In other words $\mu(\mathbf{X}; \mathbf{I})$ is the maximum normalized scalar product between distinct columns of \mathbf{X} :

$$\mu_*(\mathbf{X}; \mathbf{I}) = \frac{1}{n} \max_{i \neq j} |\langle \mathbf{X}e_i, \mathbf{X}e_j \rangle|. \quad (9)$$

The quantity (9) is known as the *coherence parameter* of the matrix \mathbf{X}/\sqrt{n} and was first defined in the context of approximation theory by Mallat and Zhang [MZ93], and by Donoho and Huo [DH01].

Assuming, for the sake of simplicity, that the columns of \mathbf{X} are normalized so that $\|\mathbf{X}e_i\|_2 = \sqrt{n}$, a small value of the coherence parameter $\mu_*(\mathbf{X}; \mathbf{I})$ means that the columns of \mathbf{X} are roughly orthogonal. We emphasize however that $\mu_*(\mathbf{X}; M)$ can be much smaller than its classical coherence parameter $\mu_*(\mathbf{X}; \mathbf{I})$. For instance, $\mu_*(\mathbf{X}; \mathbf{I}) = 0$ if and only if \mathbf{X}/\sqrt{n} is an orthogonal matrix. On the other hand, $\mu_{\min}(\mathbf{X}) = 0$ if and only if \mathbf{X} has rank² p .

The following theorem is a slight generalization of a result of [vdGBRD13]. Let us emphasize that it applies to deterministic design matrices \mathbf{X} .

Theorem 2.3. *Let $\mathbf{X} \in \mathbb{R}^{n \times p}$ be any (deterministic) design matrix, and $\hat{\theta}^* = \hat{\theta}^*(Y, \mathbf{X}; M, \lambda)$ be a general debiased estimator as per Eq. (7). Then, setting $Z = M\mathbf{X}^\top W/\sqrt{n}$, we have*

$$\sqrt{n}(\hat{\theta}^* - \theta_0) = Z + \Delta, \quad Z \sim \mathbf{N}(0, \sigma^2 M \widehat{\Sigma} M^\top), \quad \Delta = \sqrt{n}(M \widehat{\Sigma} - \mathbf{I})(\theta_0 - \hat{\theta}^n). \quad (10)$$

Further, assume that \mathbf{X} satisfies the compatibility condition for the set $S = \text{supp}(\theta_0)$, $|S| \leq s_0$, with constant ϕ_0 , and has generalized coherence parameter $\mu_* = \mu_*(\mathbf{X}; M)$, and let $K \equiv \max_{i \in [p]} (\mathbf{X}^\top \mathbf{X}/n)_{ii}$. Then, letting $\lambda = \sigma \sqrt{(c^2 \log p)/n}$, we have

$$\mathbb{P}\left(\|\Delta\|_\infty \geq \frac{4c\mu_*\sigma s_0}{\phi_0^2} \sqrt{\log p}\right) \leq 2p^{-c_0}, \quad c_0 = \frac{c^2}{32K} - 1. \quad (11)$$

Further, if $M = M_{\min}(\mathbf{X})$ minimizes the convex cost function $|M \widehat{\Sigma} - \mathbf{I}|_\infty$, then μ_* can be replaced by $\mu_{\min}(\mathbf{X})$ in Eq. (11).

The above theorem decomposes the estimation error $(\hat{\theta}^* - \theta_0)$ into a zero mean Gaussian term Z/\sqrt{n} and a bias term Δ/\sqrt{n} whose maximum entry is bounded as per Eq. (11). This estimate on $\|\Delta\|_\infty$ depends on the design matrix through two constants: the compatibility constant ϕ_0 and the generalized coherence parameter $\mu_*(\mathbf{X}; M)$. The former is a well studied property of the design matrix [BvdG11, vdGB09], and assuming ϕ_0 of order one is nearly necessary for the LASSO to achieve optimal estimation rate in high dimension. On the contrary, the definition of $\mu_*(\mathbf{X}; M)$ is a new contribution of the present paper.

The next theorem establishes that, for a natural probabilistic model of the design matrix \mathbf{X} , both ϕ_0 and $\mu_*(\mathbf{X}; M)$ can be bounded with probability converging rapidly to one as $n, p \rightarrow \infty$. Further, the bound on $\mu_*(\mathbf{X}; M)$ hold for the special choice of M that is constructed by Algorithm 1.

Theorem 2.4. *Let $\Sigma \in \mathbb{R}^{p \times p}$ be such that $\sigma_{\min}(\Sigma) \geq C_{\min} > 0$, and $\sigma_{\max}(\Sigma) \leq C_{\max} < \infty$, and $\max_{i \in [p]} \Sigma_{ii} \leq 1$. Assume $\mathbf{X}\Sigma^{-1/2}$ to have independent subgaussian rows, with zero mean and subgaussian norm $\|\Sigma^{-1/2} X_1\|_{\psi_2} = \kappa$, for some constant $\kappa \in (0, \infty)$.*

- (a) *For $\phi_0, s_0, K \in \mathbb{R}_{>0}$, let $\mathcal{E}_n = \mathcal{E}_n(\phi_0, s_0, K)$ be the event that the compatibility condition holds for $\widehat{\Sigma} = (\mathbf{X}^\top \mathbf{X}/n)$, for all sets $S \subseteq [p]$, $|S| \leq s_0$ with constant $\phi_0 > 0$, and that $\max_{i \in [p]} \widehat{\Sigma}_{i,i} \leq K$. Explicitly*

$$\mathcal{E}_n(\phi_0, s_0, K) \equiv \left\{ \mathbf{X} \in \mathbb{R}^{n \times p} : \min_{S: |S| \leq s_0} \phi(\widehat{\Sigma}, S) \geq \phi_0, \max_{i \in [p]} \widehat{\Sigma}_{i,i} \leq K, \widehat{\Sigma} = (\mathbf{X}^\top \mathbf{X}/n) \right\}. \quad (12)$$

²Of course this example requires $n \geq p$. It is the simplest example that illustrates the difference between coherence and generalized coherence, and it is not hard to find related examples with $n < p$.

Then there exists $c_* \leq 2000$ such that the following happens. If $n \geq \nu_0 s_0 \log(p/s_0)$, $\nu_0 \equiv 4c_*(C_{\max}\kappa^4/C_{\min})$, $\phi_0 = C_{\min}^{1/2}/2$, and $K \geq 1 + 20\kappa^2 \sqrt{(\log p)/n}$, then

$$\mathbb{P}(\mathbf{X} \in \mathcal{E}_n(\phi_0, s_0, K)) \geq 1 - 4e^{-c_1 n}, \quad c_1 \equiv \frac{1}{c_* \kappa^4}. \quad (13)$$

(b) For $a > 0$, $\mathcal{G}_n = \mathcal{G}_n(a)$ be the event that the problem (4) is feasible for $\mu = a\sqrt{(\log p)/n}$, or equivalently

$$\mathcal{G}_n(a) \equiv \left\{ \mathbf{X} \in \mathbb{R}^{n \times p} : \mu_{\min}(\mathbf{X}) < a\sqrt{\frac{\log p}{n}} \right\}. \quad (14)$$

Then, for $n \geq a^2 C_{\min} \log p / (4e^2 C_{\max} \kappa^4)$

$$\mathbb{P}(\mathbf{X} \in \mathcal{G}_n(a)) \geq 1 - 2p^{-c_2}, \quad c_2 \equiv \frac{a^2 C_{\min}}{24e^2 \kappa^4 C_{\max}} - 2. \quad (15)$$

The proof of this theorem is given in Section 6.2 (for part (a)) and Section 6.3 (part (b)).

The proof that event \mathcal{E}_n holds with high probability relies crucially on a theorem by Rudelson and Zhou [RZ13, Theorem 6]. Simplifying somewhat, the latter states that, if the restricted eigenvalue condition of [BRT09] holds for the population covariance Σ , then it holds with high probability for the sample covariance $\widehat{\Sigma}$. (Recall that the restricted eigenvalue condition is implied by a lower bound on the minimum singular value³, and that it implies the compatibility condition [vdGB09].)

Finally, by putting together Theorem 2.3 and Theorem 2.4, we obtain the following conclusion.

Theorem 2.5. Consider the linear model (1) and let $\widehat{\theta}^u$ be defined as per Eq. (5) in Algorithm 1, with $\mu = a\sqrt{(\log p)/n}$. Then, setting $Z = M\mathbf{X}^\top W/\sqrt{n}$, we have

$$\sqrt{n}(\widehat{\theta}^u - \theta_0) = Z + \Delta, \quad Z|\mathbf{X} \sim \mathcal{N}(0, \sigma^2 M\widehat{\Sigma}M^\top), \quad \Delta = \sqrt{n}(M\widehat{\Sigma} - \mathbf{I})(\theta_0 - \widehat{\theta}^n). \quad (16)$$

Further, under the assumptions of Theorem 2.4, and for $n \geq \max(\nu_0 s_0 \log(p/s_0), \nu_1 \log p)$, $\nu_1 = \max(1600\kappa^4, a/4)$, and $\lambda = \sigma\sqrt{(c^2 \log p)/n}$, we have

$$\mathbb{P}\left\{ \|\Delta\|_\infty \geq \left(\frac{16ac\sigma}{C_{\min}} \right) \frac{s_0 \log p}{\sqrt{n}} \right\} \leq 4e^{-c_1 n} + 4p^{-\tilde{c}_0 \wedge c_2}. \quad (17)$$

where $\tilde{c}_0 = (c^2/48) - 1$ and c_1, c_2 are given by Eqs. (13) and (15).

Finally, the tail bound (17) holds for any choice of M that is only function of the design matrix \mathbf{X} , and satisfies the feasibility condition in Eq. (4), i.e. $|M\widehat{\Sigma} - \mathbf{I}|_\infty \leq \mu$.

Assuming σ, C_{\min} of order one, the last theorem establishes that, for random designs, the maximum size of the ‘bias term’ Δ_i over $i \in [p]$ is:

$$\|\Delta\|_\infty = O\left(\frac{s_0 \log p}{\sqrt{n}} \right) \quad (18)$$

³Note, in particular, at the cost of further complicating the last statement, the condition $\sigma_{\min}(\Sigma) = \Omega(1)$ can be further weakened.

On the other hand, the ‘noise term’ Z_i is roughly of order $\sqrt{[M\widehat{\Sigma}M^\top]_{ii}}$. Bounds on the variances $[M\widehat{\Sigma}M^\top]_{ii}$ will be given in Section 3.3 showing that, if M is computed through Algorithm 1, $[M\widehat{\Sigma}M^\top]_{ii}$ is of order one for a broad family of random designs. As a consequence $|\Delta_i|$ is much smaller than $|Z_i|$ whenever $s_0 = o(\sqrt{n}/\log p)$. We summarize these remarks below.

Remark 2.6. *Theorem 2.5 only requires that the support size satisfies $s_0 = O(n/\log p)$. If we further assume $s_0 = o(\sqrt{n}/\log p)$, then we have $\|\Delta\|_\infty = o(1)$ with high probability. Hence, $\widehat{\theta}^u$ is an asymptotically unbiased estimator for θ_0 .*

A more formal comparison of the bias of $\widehat{\theta}^u$, and of the one of the LASSO estimator $\widehat{\theta}^n$ can be found in Section 2.2 below. Section 2.3 compares our approach with the related one in [vdGBRD13].

As it can be seen from the statement of Theorem 2.3 and Theorem 2.4, the claim of Theorem 2.5 does not rely on the specific choice of the objective function in optimization problem (4) and only uses the constraint on $\|\widehat{\Sigma}m - e_i\|_\infty$. In particular it holds for any matrix M that is feasible. On the other hand, the specific objective function problem (4) minimizes the variance of the noise term $\text{Var}(Z_i)$.

2.2 Discussion: The bias of the LASSO

Theorems 2.3 and 2.4 provide a quantitative framework to discuss in what sense the LASSO estimator $\widehat{\theta}^n$ is asymptotically biased, while the de-biased estimator $\widehat{\theta}^u$ is asymptotically unbiased.

Given an estimator $\widehat{\theta}^n$ of the parameter vector θ_0 , we define its *bias* to be the vector

$$\text{Bias}(\widehat{\theta}^n) \equiv \mathbb{E}\{\widehat{\theta}^n - \theta_0 | \mathbf{X}\}. \quad (19)$$

Note that, if the design is random, $\text{Bias}(\widehat{\theta}^n)$ is a measurable function of \mathbf{X} . If the design is deterministic, $\text{Bias}(\widehat{\theta}^n)$ is a deterministic quantity as well, and the conditioning is redundant.

It follows from Eq. (10) that

$$\text{Bias}(\widehat{\theta}^u) = \frac{1}{\sqrt{n}} \mathbb{E}\{\Delta | \mathbf{X}\}. \quad (20)$$

Theorem 2.5 with high probability, $\|\Delta\|_\infty = O(s_0 \log p / \sqrt{n})$. The next corollary establishes that this translates into a bound on $\text{Bias}(\widehat{\theta}^u)$ for all \mathbf{X} in a set that has probability rapidly converging to one as n, p get large.

Corollary 2.7. *Under the assumptions of Theorem 2.5, let c_1, c_2 be defined as per Eqs. (13), (15). Then we have*

$$\mathbf{X} \in \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a) \Rightarrow \|\text{Bias}(\widehat{\theta}^u)\|_\infty \leq \frac{160a}{C_{\min}} \frac{\sigma s_0 \log p}{n}, \quad (21)$$

$$\mathbb{P}\left(\mathbf{X} \in \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)\right) \geq 1 - 4e^{-n/c_*} - 2p^{-c_2}. \quad (22)$$

The proof of this corollary can be found in Appendix B.1.

This result can be contrasted with a converse result for the LASSO estimator. Namely, as stated below, there are choices of the vector θ_0 , and of the design covariance Σ , such that $\text{Bias}(\widehat{\theta}^n)$ is the sum of two terms. One is of order $\lambda = c\sigma\sqrt{(\log p)/n}$ and the second is of order $\|\text{Bias}(\widehat{\theta}^u)\|_\infty$. If

s_0 is significantly smaller than $\sqrt{n/\log p}$ (which is the main regime studied in the rest of the paper), the first term dominates and $\|\text{Bias}(\hat{\theta}^n)\|_\infty$ is much larger than $\|\text{Bias}(\hat{\theta}^u)\|_\infty$. If on the other hand s_0 is significantly larger than $\sqrt{n/\log p}$ then $\|\text{Bias}(\hat{\theta}^n)\|_\infty$ is of the same order as $\|\text{Bias}(\hat{\theta}^u)\|_\infty$. This justifies referring to $\hat{\theta}^u$ as to an *unbiased estimator*.

Notice that, since we want to establish a negative result about the LASSO, it is sufficient to exhibit a specific covariance structure Σ satisfying the assumptions of the previous corollary. Remarkably it is sufficient to consider standard designs, i.e. $\Sigma = \mathbf{I}_{p \times p}$.

Corollary 2.8. *Under the assumptions of Theorem 2.5, further consider the case $\Sigma = \mathbf{I}$. Then, there exists a numerical constant $c_{**} > 0$, a set of design matrices $\mathcal{B}_n \subseteq \mathbb{R}^{n \times p}$, and coefficient vectors $\theta_0 \in \mathbb{R}^p$, $\|\theta_0\|_0 \leq s_0$, such that*

$$\mathbf{X} \in \mathcal{B}_n \Rightarrow \|\text{Bias}(\hat{\theta}^n)\|_\infty \geq \left| \frac{2}{3}\lambda - \|\text{Bias}(\hat{\theta}^u)\|_\infty \right|, \quad (23)$$

$$\mathbb{P}(\mathcal{B}_n) \geq 1 - 6e^{-n/c_*} - 2p^{-3}. \quad (24)$$

$$(25)$$

In particular $\|\text{Bias}(\hat{\theta}^u)\|_\infty \leq \lambda/3$ (which follows from $(s_0^2 \log p)/n \leq (c/(3c_{**}))^2$) then we have

$$\|\text{Bias}(\hat{\theta}^n)\|_\infty \geq \frac{c\sigma}{3} \sqrt{\frac{\log p}{n}} \gg \|\text{Bias}(\hat{\theta}^u)\|_\infty. \quad (26)$$

On the other hand, if $\|\text{Bias}(\hat{\theta}^u)\|_\infty \geq \lambda$, then

$$\|\text{Bias}(\hat{\theta}^n)\|_\infty \geq \frac{1}{3}\|\text{Bias}(\hat{\theta}^u)\|_\infty. \quad (27)$$

A formal proof of this statement is deferred to Appendix B.2, but the underlying mathematical mechanism is quite simple and instructive. Recall that the KKT conditions for the LASSO estimator (3) read

$$\frac{1}{n} \mathbf{X}^\top (Y - \mathbf{X}\hat{\theta}^n) = \lambda v(\hat{\theta}^n), \quad (28)$$

with $v(\hat{\theta}^n) \in \mathbb{R}^p$ a vector in the subgradient of the ℓ_1 norm at $\hat{\theta}^n$. Adding $\hat{\theta}^n - \theta_0$ to both sides, and taking expectation over the noise, we get

$$\text{Bias}(\hat{\theta}^*) = \text{Bias}(\hat{\theta}^n) + \lambda \mathbb{E}\{v(\hat{\theta}^n)|\mathbf{X}\}, \quad (29)$$

Where $\hat{\theta}^*$ a debiased estimator of the general form Eq. (7), for $M = \mathbf{I}$. This suggests that $\text{Bias}(\hat{\theta}^n)$ can be decomposed in two contributions as described above, and as shown formally in Appendix B.2,

2.3 Comparison with earlier results

In this Section we briefly compare the above debiasing procedure and in particular Theorems 2.3, 2.4 and 2.5 to the results of [vdGBRD13]. In the case of linear statistical models considered here, the authors of [vdGBRD13] construct a debiased estimator of the form (7). However, instead of solving the optimization problem (4), they follow [ZZ14] and use the regression coefficients of the

i -th column of \mathbf{X} on the other columns to construct the i -th row of M . These regression coefficients are computed –once again– using the LASSO (node-wise LASSO).

It useful to spell out the most important differences between our contribution and the ones of [vdGBRD13]:

1. The case of fixed non-random designs is covered by [vdGBRD13, Theorem 2.1], which should be compared to our Theorem 2.3. While in our case the bias is controlled by the generalized coherence parameter, a similar role is played in [vdGBRD13] by the regularization parameters of the nodewise LASSO.
2. The case of random designs is covered by [vdGBRD13, Theorem 2.2, Theorem 2.4], which should be compared with our Theorem 2.5. In this case, the assumptions underlying our result are less restrictive. More precisely:
 - (a) [vdGBRD13, Theorem 2.2, Theorem 2.4] assume \mathbf{X} has i.i.d. rows, while we only assume the rows are independent.
 - (b) [vdGBRD13, Theorem 2.2, Theorem 2.4] assumes the rows of the inverse covariance matrix Σ^{-1} are sparse. More precisely, letting s_j be the number of non-zero entries of the j -th row of Σ^{-1} , [vdGBRD13] assumes $\max_{j \in [p]} s_j = o(n/\log p)$, that is much smaller than p . We do not make any sparsity assumption on Σ^{-1} , and s_j can be as large as p . [vdGBRD13, Theorem 2.4] also considers a slightly different setting, where \mathbf{X} has bounded entries, under analogous sparsity assumptions.

It is currently unknown whether the sparsity assumption in [vdGBRD13] is required for that approach to work, or it is rather an artifact of the specific analysis. Indeed [vdGBRD13, Theorem 2.1] can in principle be used to weaken this condition.

In addition our Theorem 2.5 provides the specific dependence on the maximum and minimum singular value of $\hat{\Sigma}$.

Note that solving the convex problem (4) is not more burdensome than solving the nodewise LASSO as in [ZZ14, vdGBRD13], This can be confirmed by checking that the dual of problem (4) is an ℓ_1 -regularized quadratic optimization problem. It has therefore the same complexity as the nodewise LASSO (but it is different from the nodewise LASSO).

3 Statistical inference

A direct application of Theorem 2.5 is to derive confidence intervals and statistical hypothesis tests for high-dimensional models. Throughout, we make the sparsity assumption $s_0 = o(\sqrt{n}/\log p)$ and omit explicit constants that can be readily derived from Theorem 2.5.

3.1 Preliminary lemmas

As discussed above, the bias term Δ is negligible with respect to the random term Z in the decomposition (16), provided the latter has variance of order one. Our first lemma establishes that this is indeed the case.

Lemma 3.1. Let $M = (m_1, \dots, m_p)^\top$ be the matrix with rows m_i^\top obtained by solving convex program (4) in Algorithm 1. Then for all $i \in [p]$,

$$[M\widehat{\Sigma}M^\top]_{i,i} \geq \frac{(1-\mu)^2}{\widehat{\Sigma}_{i,i}}.$$

Lemma 3.1 is proved in Appendix A.1.

Using this fact, we can then characterize the asymptotic distribution of the residuals $(\widehat{\theta}^u - \theta_{0,i})$. Theorem 2.5 naturally suggests to consider the scaled residual $\sqrt{n}(\widehat{\theta}_i^u - \theta_{0,i})/(\sigma[M\widehat{\Sigma}M^\top]_{i,i}^{1/2})$. In the next lemma we consider a slightly more general scaling, replacing σ by a consistent estimator $\widehat{\sigma}$.

Lemma 3.2. Consider a sequence of design matrices $\mathbf{X} \in \mathbb{R}^{n \times p}$, with dimensions $n \rightarrow \infty$, $p = p(n) \rightarrow \infty$ satisfying the following assumptions, for constants $C_{\min}, C_{\max}, \kappa \in (0, \infty)$ independent of n . For each n , $\Sigma \in \mathbb{R}^{p \times p}$ is such that $\sigma_{\min}(\Sigma) \geq C_{\min} > 0$, and $\sigma_{\max}(\Sigma) \leq C_{\max} < \infty$, and $\max_{i \in [p]} \Sigma_{ii} \leq 1$. Assume $\mathbf{X}\Sigma^{-1/2}$ to have independent subgaussian rows, with zero mean and subgaussian norm $\|\Sigma^{-1/2}X_1\|_{\psi_2} \leq \kappa$,

Consider the linear model (1) and let $\widehat{\theta}^u$ be defined as per Eq. (5) in Algorithm 1, with $\mu = a\sqrt{(\log p)/n}$ and $\lambda = \sigma\sqrt{(c^2 \log p)/n}$, with a, c large enough constants. Finally, let $\widehat{\sigma} = \widehat{\sigma}(y, \mathbf{X})$ an estimator of the noise level satisfying, for any $\varepsilon > 0$,

$$\lim_{n \rightarrow \infty} \sup_{\theta_0 \in \mathbb{R}^p; \|\theta_0\|_0 \leq s_0} \mathbb{P}\left(\left|\frac{\widehat{\sigma}}{\sigma} - 1\right| \geq \varepsilon\right) = 0. \quad (30)$$

If $s_0 = o(\sqrt{n}/\log p)$ ($s_0 \geq 1$), then, for all $x \in \mathbb{R}$, we have

$$\lim_{n \rightarrow \infty} \sup_{\theta_0 \in \mathbb{R}^p; \|\theta_0\|_0 \leq s_0} \left| \mathbb{P}\left\{ \frac{\sqrt{n}(\widehat{\theta}_i^u - \theta_{0,i})}{\widehat{\sigma}[M\widehat{\Sigma}M^\top]_{i,i}^{1/2}} \leq x \right\} - \Phi(x) \right| = 0. \quad (31)$$

The proof of this lemma can be found in Section 6.5. We also note that the dependence of a, c on $C_{\min}, C_{\max}, \kappa$ can be easily reconstructed from Theorem 2.4.

The last lemma requires a consistent estimator of σ , in the sense of Eq. (30). Several proposal have been made to estimate the noise level in high-dimensional linear regression. A short list of references includes [FL01, FL08, SBvdG10, Zha10, SZ12, BC13, FGH12, RTF13, Dic12, FSW09, BEM13]. Consistency results have been proved or can be proved for several of these estimators.

In order to demonstrate that the consistency criterion (30) can be achieved, we use the scaled LASSO [SZ12] given by

$$\{\widehat{\theta}^n(\widetilde{\lambda}), \widehat{\sigma}(\widetilde{\lambda})\} \equiv \arg \min_{\theta \in \mathbb{R}^p, \sigma > 0} \left\{ \frac{1}{2\sigma n} \|Y - \mathbf{X}\theta\|_2^2 + \frac{\sigma}{2} + \widetilde{\lambda} \|\theta\|_1 \right\}. \quad (32)$$

This is a joint convex optimization problem which provides an estimate of the noise level in addition to an estimate of θ_0 .

The following lemma uses the analysis of [SZ12] to show that $\widehat{\sigma}$ thus defined satisfies the consistency criterion (30).

Lemma 3.3. Under the assumptions of Lemma 3.2, let $\widehat{\sigma} = \widehat{\sigma}(\widetilde{\lambda})$ be the scaled LASSO estimator of the noise level, see Eq. (32), with $\widetilde{\lambda} = 10\sqrt{(2 \log p)/n}$. Then $\widehat{\sigma}$ thus satisfies Eq. (30).

The proof of this lemma is fairly straightforward and can be found in Appendix C.

3.2 Confidence intervals

In view of Lemma 3.2, it is quite straightforward to construct asymptotically valid confidence intervals. Namely, for $i \in [p]$ and significance level $\alpha \in (0, 1)$, we let

$$\begin{aligned} J_i(\alpha) &\equiv [\hat{\theta}_i^u - \delta(\alpha, n), \hat{\theta}_i^u + \delta(\alpha, n)], \\ \delta(\alpha, n) &\equiv \Phi^{-1}(1 - \alpha/2) \frac{\hat{\sigma}}{\sqrt{n}} [M\hat{\Sigma}M^\top]_{i,i}^{1/2}. \end{aligned} \quad (33)$$

Theorem 3.4. *Consider a sequence of design matrices $\mathbf{X} \in \mathbb{R}^{n \times p}$, with dimensions $n \rightarrow \infty$, $p = p(n) \rightarrow \infty$ satisfying the assumptions of Lemma 3.2.*

Consider the linear model (1) and let $\hat{\theta}^u$ be defined as per Eq. (5) in Algorithm 1, with $\mu = a\sqrt{(\log p)/n}$ and $\lambda = \sigma\sqrt{(c^2 \log p)/n}$, with a, c large enough constants. Finally, let $\hat{\sigma} = \hat{\sigma}(y, \mathbf{X})$ a consistent estimator of the noise level in the sense of Eq. (30). Then the confidence interval $J_i(\alpha)$ is asymptotically valid, namely

$$\lim_{n \rightarrow \infty} \mathbb{P}(\theta_{0,i} \in J_i(\alpha)) = 1 - \alpha. \quad (34)$$

Proof. The proof is an immediate consequence of Lemma 3.2 since

$$\lim_{n \rightarrow \infty} \mathbb{P}(\theta_{0,i} \in J_i(\alpha)) = \lim_{n \rightarrow \infty} \mathbb{P} \left\{ \frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \leq \Phi^{-1}(1 - \alpha/2) \right\} \quad (35)$$

$$- \lim_{n \rightarrow \infty} \mathbb{P} \left\{ \frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \leq -\Phi^{-1}(1 - \alpha/2) \right\} \quad (36)$$

$$= 1 - \alpha. \quad (37)$$

□

3.3 Hypothesis testing

An important advantage of sparse linear regression models is that they provide parsimonious explanations of the data in terms of a small number of covariates. The easiest way to select the ‘active’ covariates is to choose the indexes i for which $\hat{\theta}_i^n \neq 0$. This approach however does not provide a measure of statistical significance for the finding that the coefficient is non-zero.

More precisely, we are interested in testing an individual null hypothesis $H_{0,i} : \theta_{0,i} = 0$ versus the alternative $H_{A,i} : \theta_{0,i} \neq 0$, and assigning p -values for these tests. We construct a p -value P_i for the test $H_{0,i}$ as follows:

$$P_i = 2 \left(1 - \Phi \left(\frac{\sqrt{n} |\hat{\theta}_i^u|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \right) \right). \quad (38)$$

The decision rule is then based on the p -value P_i :

$$\hat{T}_{i,\mathbf{X}}(y) = \begin{cases} 1 & \text{if } P_i \leq \alpha \quad (\text{reject } H_{0,i}), \\ 0 & \text{otherwise} \quad (\text{accept } H_{0,i}), \end{cases} \quad (39)$$

where α is the fixed target Type I error probability. We measure the quality of the test $\widehat{T}_{i,\mathbf{X}}(y)$ in terms of its significance level α_i and statistical power $1 - \beta_i$. Here α_i is the probability of type I error (i.e. of a false positive at i) and β_i is the probability of type II error (i.e. of a false negative at i).

Note that it is important to consider the tradeoff between statistical significance and power. Indeed any significance level α can be achieved by randomly rejecting $H_{0,i}$ with probability α . This test achieves power $1 - \beta = \alpha$. Further note that, without further assumption, no nontrivial power can be achieved. In fact, choosing $\theta_{0,i} \neq 0$ arbitrarily close to zero, $H_{0,i}$ becomes indistinguishable from its alternative. We will therefore assume that, whenever $\theta_{0,i} \neq 0$, we have $|\theta_{0,i}| > \gamma$ as well. We take a minimax perspective and require the test to behave uniformly well over s_0 -sparse vectors. Formally, given a family of tests $T_{i,\mathbf{X}} : \mathbb{R}^n \rightarrow \{0, 1\}$, indexed by $i \in [p]$, $\mathbf{X} \in \mathbb{R}^{n \times p}$, we define, for $\gamma > 0$ a lower bound on the non-zero entries:

$$\alpha_{i,n}(T) \equiv \sup \left\{ \mathbb{P}_{\theta_0}(T_{i,\mathbf{X}}(y) = 1) : \theta_0 \in \mathbb{R}^p, \|\theta_0\|_0 \leq s_0(n), \theta_{0,i} = 0 \right\}. \quad (40)$$

$$\beta_{i,n}(T; \gamma) \equiv \sup \left\{ \mathbb{P}_{\theta_0}(T_{i,\mathbf{X}}(y) = 0) : \theta_0 \in \mathbb{R}^p, \|\theta_0\|_0 \leq s_0(n), |\theta_{0,i}| \geq \gamma \right\}. \quad (41)$$

Here, we made dependence on n explicit. Also, $\mathbb{P}_{\theta}(\cdot)$ denotes the induced probability for random design \mathbf{X} and noise realization w , given the fixed parameter vector θ . Our next theorem establishes bounds on $\alpha_{i,n}(\widehat{T})$ and $\beta_{i,n}(\widehat{T}; \gamma)$ for our decision rule (39).

Theorem 3.5. *Consider a sequence of design matrices $\mathbf{X} \in \mathbb{R}^{n \times p}$, with dimensions $n \rightarrow \infty$, $p = p(n) \rightarrow \infty$ satisfying the assumptions of Lemma 3.2.*

Consider the linear model (1) and let $\widehat{\theta}^u$ be defined as per Eq. (5) in Algorithm 1, with $\mu = a\sqrt{(\log p)/n}$ and $\lambda = \sigma\sqrt{(c^2 \log p)/n}$, with a, c large enough constants. Finally, let $\widehat{\sigma} = \widehat{\sigma}(y, \mathbf{X})$ a consistent estimator of the noise level in the sense of Eq. (30), and \widehat{T} be the test defined in Eq. (39).

Then the following holds true for any fixed sequence of integers $i = i(n)$:

$$\lim_{n \rightarrow \infty} \alpha_{i,n}(\widehat{T}) \leq \alpha. \quad (42)$$

$$\liminf_{n \rightarrow \infty} \frac{1 - \beta_{i,n}(\widehat{T}; \gamma)}{1 - \beta_{i,n}^*(\gamma)} \geq 1, \quad 1 - \beta_{i,n}^*(\gamma) \equiv G\left(\alpha, \frac{\sqrt{n}\gamma}{\sigma_{[\Sigma_{i,i}^{-1}]^{1/2}}}\right), \quad (43)$$

where, for $\alpha \in [0, 1]$ and $u \in \mathbb{R}_+$, the function $G(\alpha, u)$ is defined as follows:

$$G(\alpha, u) = 2 - \Phi\left(\Phi^{-1}\left(1 - \frac{\alpha}{2}\right) + u\right) - \Phi\left(\Phi^{-1}\left(1 - \frac{\alpha}{2}\right) - u\right).$$

Theorem 3.5 is proved in Appendix 6.6. It is easy to see that, for any $\alpha > 0$, $u \mapsto G(\alpha, u)$ is continuous and monotone increasing. Moreover, $G(\alpha, 0) = \alpha$ which is the trivial power obtained by randomly rejecting $H_{0,i}$ with probability α . As γ deviates from zero, we obtain nontrivial power. Notice that in order to achieve a specific power $\beta > \alpha$, our scheme requires $\gamma \geq c_\beta(\sigma/\sqrt{n})$, for some constant c_β that depends on β . This is because $\Sigma_{i,i}^{-1} \leq \sigma_{\max}(\Sigma^{-1}) \leq (\sigma_{\min}(\Sigma))^{-1} = O(1)$.

3.3.1 Near optimality of the hypothesis testing procedure

The authors of [JM13b] prove an upper bound for the minimax power of tests with a given significance level α , under random designs. For the readers' convenience, we recall here this result. (The following is a restatement of [JM13b, Theorem 2.3], together with a standard estimate on the tail of chi-squared random variables.)

Theorem 3.6 ([JM13b]). Assume $\mathbf{X} \in \mathbb{R}^{n \times p}$ to be a random design matrix with i.i.d. Gaussian rows with zero mean and covariance Σ . For $i \in [p]$, let $T_{i,\mathbf{X}} : \mathbb{R}^n \rightarrow \mathbb{R}^n$ be a hypothesis testing procedure for testing $H_{0,i} : \theta_{0,i} = 0$, and denote by $\alpha_i(T)$ and $\beta_{i,n}(T; \gamma)$ its fraction of type I and type II errors, cf. Eqs. (40) and (41). Finally, for $S \subseteq [p] \setminus \{i\}$, define $\Sigma_{i|S} \equiv \Sigma_{ii} - \Sigma_{i,S} \Sigma_{S,S}^{-1} \Sigma_{S,i} \in \mathbb{R}$.

For any $\ell \in \mathbb{R}$ and $|S| < s_0 < n$, if $\alpha_{i,n}(T) \leq \alpha$, then

$$1 - \beta_{i,n}(T; \gamma) \leq G\left(\alpha, \frac{\gamma}{\sigma_{\text{eff}}(\xi)}\right) + e^{-\xi^2/8}, \quad (44)$$

$$\sigma_{\text{eff}}(\xi) \equiv \frac{\sigma}{\Sigma_{i|S}^{1/2}(\sqrt{n - s_0 + 1} + \xi)}, \quad (45)$$

for any $\xi \in [0, (3/2)\sqrt{n - s_0 + 1}]$.

The intuition behind this bound is straightforward: the power of any test for $H_{0,i} : \theta_{0,i} = 0$ is upper bounded by the power of an oracle test that is given access to the support of θ_0 , with the eventual exclusion of i . Namely, the oracle has access to $\text{supp}(\theta_0) \setminus \{i\}$ and outputs a test for $H_{0,i}$. Computing the minimax power of such oracle reduces to a classical hypothesis testing problem.

Let us emphasize that the last theorem applies to *Gaussian* random designs. Since this theorem establishes a negative result (an upper bound on power) it makes sense to consider this somewhat more specialized setting.

Using this upper bound, we can restate Theorem 3.5 as follows.

Corollary 3.7. Consider a Gaussian random design model that satisfies the conditions of Theorem 3.5, and let \hat{T} be the testing procedure defined in Eq. (39), with $\hat{\theta}^u$ as in Algorithm 1. Further, let

$$\eta_{\Sigma, s_0} \equiv \min_{i \in [p]; S} \left\{ \Sigma_{i|S} \Sigma_{ii}^{-1} : S \subseteq [p] \setminus \{i\}, |S| < s_0 \right\}. \quad (46)$$

Under the sparsity assumption $s_0 = o(\sqrt{n}/\log p)$, the following holds true. If $\{T_{i,\mathbf{X}}\}$ is any sequence of tests with $\limsup_{n \rightarrow \infty} \alpha_{i,n}(T) \leq \alpha$, then

$$\liminf_{n \rightarrow \infty} \frac{1 - \beta_{i,n}(\hat{T}; \gamma)}{1 - \beta_{i,n/\eta_{\Sigma, s_0}}(T; \gamma)} \geq 1. \quad (47)$$

In other words, the asymptotic efficiency of the test \hat{T} is at least $1/\eta_{\Sigma, s_0}$.

Hence, our test \hat{T} has nearly optimal power in the following sense. It has power at least as large as the power of any other test T , provided the latter is applied to a sample size increased by a factor η_{Σ, s_0} .

Further, under the assumptions of Theorem 2.5, the factor η_{Σ, s_0} is a bounded constant. Indeed

$$\eta_{\Sigma, s_0} \leq \Sigma_{i,i}^{-1} \Sigma_{i,i} \leq \frac{\sigma_{\max}(\Sigma)}{\sigma_{\min}(\Sigma)} \leq \frac{C_{\max}}{C_{\min}}, \quad (48)$$

since $\Sigma_{ii}^{-1} \leq (\sigma_{\min}(\Sigma))^{-1}$, and $\Sigma_{i|S} \leq \Sigma_{i,i} \leq \sigma_{\max}(\Sigma)$ due to $\Sigma_{S,S} \succ 0$.

Note that n , γ and σ appears in our upper bound (44) in the combination $\gamma\sqrt{n}/\sigma$, which is the natural measure of the signal-to-noise ratio (where, for simplicity, we neglected $s_0 = o(\sqrt{n}/\log p)$ with respect to n). Hence, the above result can be restated as follows. The test \hat{T} has power at least as large as the power of any other test T , provided the latter is applied at a noise level augmented by a factor $\sqrt{\eta_{\Sigma, s_0}}$.

3.4 Generalization to simultaneous confidence intervals

In many situations, it is necessary to perform statistical inference on more than one of the parameters simultaneously. For instance, we might be interested in performing inference about $\theta_{0,R} \equiv (\theta_{0,i})_{i \in R}$ for some set $R \subseteq [p]$.

The simplest generalization of our method is to the case in which $|R|$ stays finite as $n, p \rightarrow \infty$. In this case we have the following generalization of Lemma 3.2. (The proof is the same as for Lemma 3.2, and hence we omit it.)

Lemma 3.8. *Under the assumptions of Lemma 3.2, define*

$$Q^{(n)} \equiv \frac{\hat{\sigma}^2}{n} [M\widehat{\Sigma}M^\top]. \quad (49)$$

Let $R = R(n)$ be a sequence of sets $R(n) \subseteq [p]$, with $|R(n)| = k$ fixed as $n, p \rightarrow \infty$, and further assume $s_0 = o(\sqrt{n}/\log p)$, with $s_0 \geq 1$. Then, for all $x = (x_1, \dots, x_k) \in \mathbb{R}^k$, we have

$$\lim_{n \rightarrow \infty} \sup_{\theta_0 \in \mathbb{R}^p; \|\theta_0\|_0 \leq s_0} \left| \mathbb{P} \left\{ (Q_{R,R}^{(n)})^{-1/2} (\hat{\theta}_R^u - \theta_{0,R}) \leq x \right\} - \Phi_k(x) \right| = 0, \quad (50)$$

where $(a_1, \dots, a_k) \leq (b_1, \dots, b_k)$ indicates that $a_1 \leq b_1, \dots, a_k \leq b_k$, and $\Phi_k(x) = \Phi(x_1) \cdots \Phi(x_k)$.

This lemma allows to construct confidence regions for low-dimensional projections of θ_0 , much in the same way as we used Lemma 3.2 to compute confidence intervals for one-dimensional projections in Section 3.2.

Explicitly, let $\mathcal{C}_{k,\alpha} \subseteq \mathbb{R}^k$ be any Borel set such that $\int_{\mathcal{C}_{k,\alpha}} \phi_k(x) dx \geq 1 - \alpha$, where

$$\phi_k(x) = \frac{1}{(2\pi)^{k/2}} \exp\left(-\frac{\|x\|^2}{2}\right),$$

is the k -dimensional Gaussian density. Then, for $R \subseteq [p]$, we define $J_R(\alpha) \subseteq \mathbb{R}^k$ as follows

$$J_R(\alpha) \equiv \hat{\theta}_R^u + (Q_{R,R}^{(n)})^{1/2} \mathcal{C}_{k,\alpha}. \quad (51)$$

Then Lemma 3.8 implies (under the assumptions stated there) that $J_R(\alpha)$ is a valid confidence region

$$\lim_{n \rightarrow \infty} \mathbb{P}(\theta_{0,R} \in J_R(\alpha)) = 1 - \alpha. \quad (52)$$

A more challenging regime is the one of large-scale inference, that corresponds to $|R(n)| \rightarrow \infty$ with n . Even in the seemingly simple case in which a correct p -value is given for each individual coordinate, the problem of aggregating them has attracted considerable amount of work, see e.g. [Efr10] for an overview.

Here we limit ourself to designing a testing procedure for the family of hypotheses $\{H_{0,i} : \theta_{0,i} = 0\}_{i \in [p]}$ that controls the familywise error rate (FWER). Namely we want to define $T_{i,\mathbf{X}} : \mathbb{R}^n \rightarrow \{0, 1\}$, for each $i \in [p]$, $\mathbf{X} \in \mathbb{R}^{n \times p}$ such that

$$\text{FWER}(T, n) \equiv \sup_{\theta_0 \in \mathbb{R}^p, \|\theta_0\|_0 \leq s_0} \mathbb{P} \left\{ \exists i \in [p] : \theta_{0,i} = 0, T_{i,\mathbf{X}}(y) = 1 \right\}, \quad (53)$$

In order to achieve familywise error control, we adopt a standard trick based on Bonferroni inequality. Given p -values defined as per Eq. (38), we let

$$\widehat{T}_{i,\mathbf{X}}^{\text{F}}(y) = \begin{cases} 1 & \text{if } P_i \leq \alpha/p & (\text{reject } H_{0,i}), \\ 0 & \text{otherwise} & (\text{accept } H_{0,i}). \end{cases} \quad (54)$$

Then we have the following error control guarantee.

Theorem 3.9. *Consider a sequence of design matrices $\mathbf{X} \in \mathbb{R}^{n \times p}$, with dimensions $n \rightarrow \infty$, $p = p(n) \rightarrow \infty$ satisfying the assumptions of Lemma 3.2.*

Consider the linear model (1) and let $\widehat{\theta}^u$ be defined as per Eq. (5) in Algorithm 1, with $\mu = a\sqrt{(\log p)/n}$ and $\lambda = \sigma\sqrt{(c^2 \log p)/n}$, with a, c large enough constants. Finally, let $\widehat{\sigma} = \widehat{\sigma}(y, \mathbf{X})$ be a consistent estimator of the noise level in the sense of Eq. (30), and \widehat{T} be the test defined in Eq. (54). Then:

$$\limsup_{n \rightarrow \infty} \text{FWER}(\widehat{T}^{\text{F}}, n) \leq \alpha. \quad (55)$$

The proof of this theorem is similar to the one of Lemma 3.2 and Theorem 3.5, and is deferred to Appendix D.

4 Non-Gaussian noise

As can be seen from the proof of Theorem 2.5, $Z = M\mathbf{X}^{\text{T}}W/\sqrt{n}$, and since the noise is Gaussian, i.e., $W \sim \mathbf{N}(0, \sigma^2\mathbf{I})$, we have $Z|\mathbf{X} \sim \mathbf{N}(0, \sigma^2 M\widehat{\Sigma}M^{\text{T}})$. We claim that the distribution of the coordinates of Z is asymptotically Gaussian, even if W is non-Gaussian, provided the definition of M is modified slightly. As a consequence, the definition of confidence intervals and p -values in Corollary 3.4 and (38) remain valid in this broader setting.

In case of non-Gaussian noise, we write

$$\begin{aligned} \frac{\sqrt{n}(\widehat{\theta}_i^u - \theta_{0,i})}{\sigma[M\widehat{\Sigma}M^{\text{T}}]_{i,i}^{1/2}} &= \frac{1}{\sqrt{n}} \frac{m_i^{\text{T}}\mathbf{X}^{\text{T}}W}{\sigma[m_i^{\text{T}}\widehat{\Sigma}m_i]^{1/2}} + o(1) \\ &= \frac{1}{\sqrt{n}} \sum_{j=1}^n \frac{m_i^{\text{T}}X_jW_j}{\sigma[m_i^{\text{T}}\widehat{\Sigma}m_i]^{1/2}} + o(1). \end{aligned}$$

Conditional on \mathbf{X} , the summands $\xi_j = m_i^{\text{T}}X_jW_j/(\sigma[m_i^{\text{T}}\widehat{\Sigma}m_i]^{1/2})$ are independent and zero mean. Further, $\sum_{j=1}^n \mathbb{E}(\xi_j^2|\mathbf{X}) = 1$. Therefore, if Lindenberg condition holds, namely for every $\varepsilon > 0$, almost surely

$$\lim_{n \rightarrow \infty} \frac{1}{n} \sum_{j=1}^n \mathbb{E}(\xi_j^2 \mathbb{I}_{\{|\xi_j| > \varepsilon\sqrt{n}\}}|\mathbf{X}) = 0, \quad (56)$$

then $\sum_{j=1}^n \xi_j/\sqrt{n}|\mathbf{X} \xrightarrow{\text{d}} \mathbf{N}(0, 1)$, from which we can build the valid p -values as in (38).

In order to ensure that the Lindeberg condition holds, we modify the optimization problem (57) as follows:

$$\begin{aligned}
& \text{minimize} && m^\top \widehat{\Sigma} m \\
& \text{subject to} && \|\widehat{\Sigma} m - e_i\|_\infty \leq \mu \\
& && \|\mathbf{X}m\|_\infty \leq n^\beta \quad \text{for arbitrary fixed } 0 < \beta < 1/2
\end{aligned} \tag{57}$$

Next theorem shows the validity of the proposed p -values in the non-Gaussian noise setting.

Theorem 4.1. *Suppose that the noise variables W_i are independent with $\mathbb{E}(W_i) = 0$, $\mathbb{E}(W_i^2) = \sigma^2$, and $\mathbb{E}(|W_i|^{2+a}) \leq C \sigma^{2+a}$ for some $a > (1/2 - \beta)^{-1}$.*

Let $M = (m_1, \dots, m_p)^\top$ be the matrix with rows m_i^\top obtained by solving optimization problem (57). Then under the assumptions of Theorem 2.5, and for sparsity level $s_0 = o(\sqrt{n}/\log p)$, an asymptotic two-sided confidence interval for $\theta_{0,i}$ with significance α is given by $I_i = [\widehat{\theta}_i^u - \delta(\alpha, n), \widehat{\theta}_i^u + \delta(\alpha, n)]$ where

$$\delta(\alpha, n) = \Phi^{-1}(1 - \alpha/2) \widehat{\sigma} n^{-1/2} \sqrt{[M \widehat{\Sigma} M^\top]_{i,i}}. \tag{58}$$

Further, an asymptotically valid p -value P_i for testing null hypothesis $H_{0,i}$ is constructed as:

$$P_i = 2 \left(1 - \Phi \left(\frac{\sqrt{n} |\widehat{\theta}_i^u|}{[M \widehat{\Sigma} M^\top]_{i,i}^{1/2}} \right) \right).$$

Theorem 4.1 is proved in Section 6.7.

5 Numerical experiments

5.1 Synthetic data

We consider linear model (2), where the rows of design matrix \mathbf{X} are fixed i.i.d. realizations from $\mathbf{N}(0, \Sigma)$, where $\Sigma \in \mathbb{R}^{p \times p}$ is a circulant symmetric matrix with entries Σ_{jk} given as follows for $j \leq k$:

$$\Sigma_{jk} = \begin{cases} 1 & \text{if } k = j, \\ 0.1 & \text{if } k \in \{j+1, \dots, j+5\} \\ & \text{or } k \in \{j+p-5, \dots, j+p-1\}, \\ 0 & \text{for all other } j \leq k. \end{cases} \tag{59}$$

Regarding the regression coefficient, we consider a uniformly random support $S \subseteq [p]$, with $|S| = s_0$ and let $\theta_{0,i} = b$ for $i \in S$ and $\theta_{0,i} = 0$ otherwise. The measurement errors are $W_i \sim \mathbf{N}(0, 1)$, for $i \in [n]$. We consider several configurations of (n, p, s_0, b) and for each configuration report our results based on 20 independent realizations of the model with fixed design and fixed regression coefficients. In other words, we repeat experiments over 20 independent realization of the measurement errors.

We use the regularization parameter $\lambda = 4\widehat{\sigma} \sqrt{(2 \log p)/n}$, where $\widehat{\sigma}$ is given by the scaled LASSO as per equation (32) with $\widetilde{\lambda} = 10\sqrt{(2 \log p)/n}$. Furthermore, parameter μ (cf. Eq. (4)) is set to

$$\mu = 2 \sqrt{\frac{\log p}{n}}.$$

This choice of μ is guided by Theorem 2.4 (b).

Throughout, we set the significance level $\alpha = 0.05$.

Confidence intervals. For each configuration, we consider 20 independent realizations of measurement noise and for each parameter $\theta_{0,i}$, we compute the average length of the corresponding confidence interval, denoted by $\text{Avglength}(J_i(\alpha))$ where $J_i(\alpha)$ is given by equation (33) and the average is taken over the realizations. We then define

$$\ell \equiv p^{-1} \sum_{i \in [p]} \text{Avglength}(J_i(\alpha)). \quad (60)$$

We also consider the average length of intervals for the active and inactive parameters, as follows:

$$\ell_S \equiv s_0^{-1} \sum_{i \in S} \text{Avglength}(J_i(\alpha)), \quad \ell_{S^c} \equiv (p - s_0)^{-1} \sum_{i \in S^c} \text{Avglength}(J_i(\alpha)). \quad (61)$$

Similarly, we consider average coverage for individual parameters. We define the following three metrics:

$$\widehat{\text{Cov}} \equiv p^{-1} \sum_{i \in [p]} \widehat{\mathbb{P}}[\theta_{0,i} \in J_i(\alpha)], \quad (62)$$

$$\widehat{\text{Cov}}_S \equiv s_0^{-1} \sum_{i \in S} \widehat{\mathbb{P}}[\theta_{0,i} \in J_i(\alpha)], \quad (63)$$

$$\widehat{\text{Cov}}_{S^c} \equiv (p - s_0)^{-1} \sum_{i \in S^c} \widehat{\mathbb{P}}[0 \in J_i(\alpha)], \quad (64)$$

where $\widehat{\mathbb{P}}$ denotes the empirical probability computed based on the 20 realizations for each configuration. The results are reported in Table 1. In Fig. 1, we plot the constructed 95%-confidence intervals for one realization of configuration $(n, p, s_0, b) = (1000, 600, 10, 1)$. For sake of clarity, we plot the confidence intervals for only 100 of the 1000 parameters.

False positive rates and statistical powers. Table 2 summarizes the false positive rates and the statistical powers achieved by our proposed method, the multisample-splitting method [MMB09], and the ridge-type projection estimator [Büh13] for several configurations. The results are obtained by taking average over 20 independent realizations of measurement errors for each configuration. As we see the multisample-splitting achieves false positive rate 0 on all of the configurations considered here, making no type I error. However, the true positive rate is always smaller than that of our proposed method. By contrast, our method achieves false positive rate close to the pre-assigned significance level $\alpha = 0.05$ and obtains much higher true positive rate. Similar to the multisample-splitting, the ridge-type projection estimator is conservative and achieves false positive rate smaller than α . This, however, comes at the cost of a smaller true positive rate than our method. It is worth noting that an ideal testing procedure should allow to control the level of statistical significance α , and obtain the maximum true positive rate at that level.

Here, we used the R-package `hdi` to test multisample-splitting and the ridge-type projection estimator.

Let $Z = (z_i)_{i=1}^p$ denote the vector with $z_i \equiv \sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})/\hat{\sigma}\sqrt{[M\widehat{\Sigma}M^\top]_{i,i}}$. Fig. 2 shows the sample quantiles of Z versus the quantiles of the standard normal distribution for one realization of the configuration $(n, p, s_0, b) = (1000, 600, 10, 1)$. The scattered points are close to the line with unit

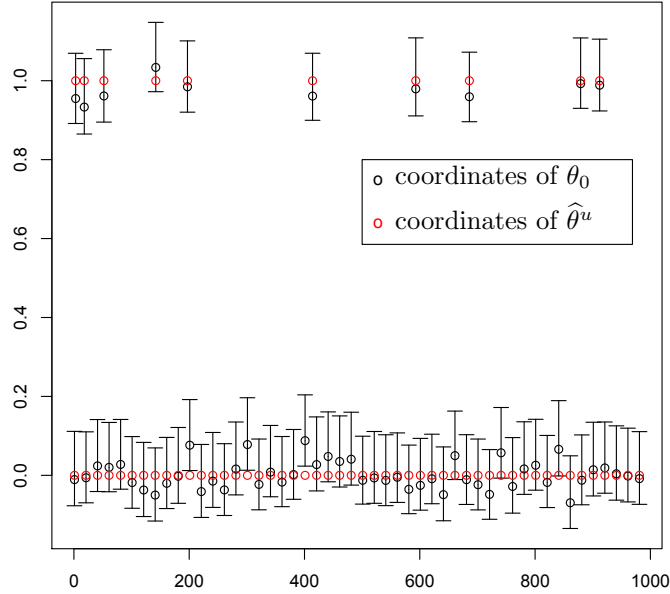


Figure 1: 95% confidence intervals for one realization of configuration $(n, p, s_0, b) = (1000, 600, 10, 1)$. For clarity, we plot the confidence intervals for only 100 of the 1000 parameters. The true parameters $\theta_{0,i}$ are in red and the coordinates of the debiased estimator $\hat{\theta}^u$ are in black.

Configuration \ Measure	Measure					
	ℓ	ℓ_S	ℓ_{Sc}	$\widehat{\text{Cov}}$	$\widehat{\text{Cov}}_S$	$\widehat{\text{Cov}}_{Sc}$
(1000, 600, 10, 0.5)	0.1870	0.1834	0.1870	0.9766	0.9600	0.9767
(1000, 600, 10, 0.25)	0.1757	0.1780	0.1757	0.9810	0.9000	0.9818
(1000, 600, 10, 0.1)	0.1809	0.1823	0.1809	0.9760	1	0.9757
(1000, 600, 30, 0.5)	0.2107	0.2108	0.2107	0.9780	0.9866	0.9777
(1000, 600, 30, 0.25)	0.1956	0.1961	0.1956	0.9660	0.9660	0.9659
(1000, 600, 30, 0.1)	0.2023	0.2043	0.2023	0.9720	0.9333	0.9732
(2000, 1500, 50, 0.5)	0.1383	0.1391	0.1383	0.9754	0.9800	0.9752
(2000, 1500, 50, 0.25)	0.1356	0.1363	0.1355	0.9720	0.9600	0.9723
(2000, 1500, 50, 0.1)	0.1361	0.1361	0.1361	0.9805	1	0.9800
(2000, 1500, 25, 0.5)	0.1233	0.1233	0.1233	0.9731	0.9680	0.9731
(2000, 1500, 25, 0.25)	0.1208	0.1208	0.1208	0.9735	1	0.9731
(2000, 1500, 25, 0.1)	0.1242	0.1237	0.1242	0.9670	0.9200	0.9676

Table 1: Simulation results for the synthetic data described in Section 5.1. The results corresponds to 95% confidence intervals.

slope and zero intercept. This confirms the result of Theorem 3.2 regarding the gaussianity of the entries z_i .

For the same problem, in Fig. 3 we plot the empirical CDF of the computed p -values restricted to the variables outside the support. Clearly, the p -values for these entries are uniformly distributed as expected.

Configuration	Our method		Multisample-splitting		Ridge-type projection estimator	
	FP	TP	FP	TP	FP	TP
(1000, 600, 10, 0.5)	0.0452	1	0	1	0.0284	0.8531
(1000, 600, 10, 0.25)	0.0393	1	0	0.4	0.02691	0.7506
(1000, 600, 10, 0.1)	0.0383	0.8	0	0	0.2638	0.6523
(1000, 600, 30, 0.5)	0.0433	1	0	1	0.0263	0.8700
(1000, 600, 30, 0.25)	0.0525	1	0	0.4	0.2844	0.8403
(1000, 600, 30, 0.1)	0.0402	0.7330	0	0	0.2238	0.6180
(2000, 1500, 50, 0.5)	0.0421	1	0	1	0.0301	0.9013
(2000, 1500, 50, 0.25)	0.0415	1	0	1	0.0292	0.8835
(2000, 1500, 50, 0.1)	0.0384	0.9400	0	0	0.02655	0.7603
(2000, 1500, 25, 0.5)	0.0509	1	0	1	0.0361	0.9101
(2000, 1500, 25, 0.25)	0.0481	1	0	1	0.3470	0.8904
(2000, 1500, 25, 0.1)	0.0551	1	0	0.16	0.0401	0.8203

Table 2: Simulation results for the synthetic data described in Section 5.1. The false positive rates (FP) and the true positive rates (TP) are computed at significance level $\alpha = 0.05$.

5.2 Real data

As a real data example, we consider a high-throughput genomic data set concerning riboflavin (vitamin B_2) production rate. This data set is made publicly available by [BKM14] and contains $n = 71$ samples and $p = 4,088$ covariates corresponding to $p = 4,088$ genes. For each sample, there is a real-valued response variable indicating the logarithm of the riboflavin production rate along with the logarithm of the expression level of the $p = 4,088$ genes as the covariates.

Following [BKM14], we model the riboflavin production rate as a linear model with $p = 4,088$ covariates and $n = 71$ samples, as in Eq. (1). We use the R package `glmnet` [FHT10] to fit the LASSO estimator. Similar to the previous section, we use the regularization parameter $\lambda = 4\hat{\sigma}\sqrt{(2\log p)/n}$, where $\hat{\sigma}$ is given by the scaled LASSO as per equation (32) with $\tilde{\lambda} = 10\sqrt{(2\log p)/n}$. This leads to the choice $\lambda = 0.036$. The resulting model contains 30 genes (plus an intercept term) corresponding to the nonzero parameters of the lasso estimator.

We use Eq. (38) to construct p -values for different genes. Adjusting FWER to 5% significance level, we find two significant genes, namely genes YXLD-at and YXLE-at. By contrast, the multisample-splitting method proposed in [MMB09] finds only the gene YXLD-at at the FWER-adjusted 5% significance level. Also the Ridge-type projection estimator, proposed in [Büh13], returns no significance gene. (See [BKM14] for further discussion on these methods.) This indicates that these methods are more conservative and produce typically larger p -values.

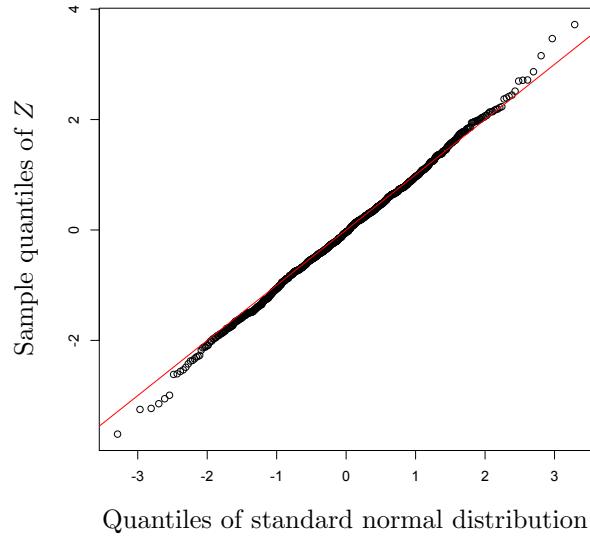


Figure 2: Q-Q plot of Z for one realization of configuration $(n, p, s_0, b) = (1000, 600, 10, 1)$.

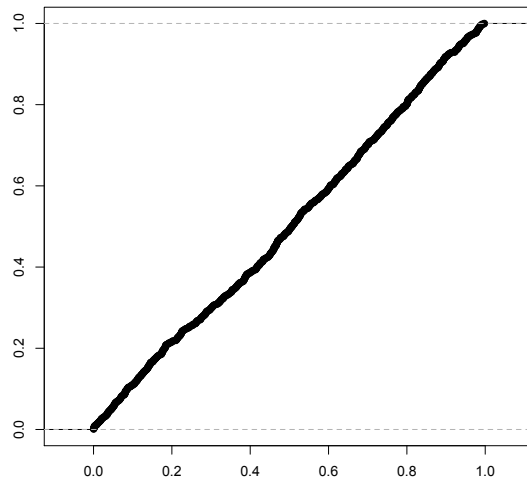


Figure 3: Empirical CDF of the computed p -values (restricted to entries outside the support) for one realization of configuration $(n, p, s_0, b) = (1000, 600, 10, 1)$. Clearly the plot confirms that the p -values are distributed according to uniform distribution.

In Fig. 4 we plot the empirical CDF of the computed p -values for riboflavin example. Clearly the plot confirms that the p -values are distributed according to uniform distribution.

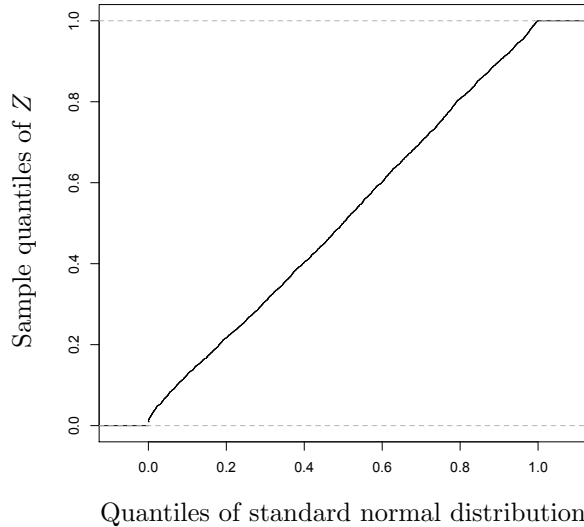


Figure 4: Empirical CDF of the computed p -values for riboflavin example. Clearly the plot confirms that the p -values are distributed according to uniform distribution.

6 Proofs

6.1 Proof of Theorem 2.3

Substituting $Y = \mathbf{X}\theta_0 + W$ in the definition (7), we get

$$\hat{\theta}^* = \hat{\theta}^n + \frac{1}{n} M \mathbf{X}^\top \mathbf{X} (\theta_0 - \hat{\theta}^n) + \frac{1}{n} M \mathbf{X}^\top W \quad (65)$$

$$= \theta_0 + \frac{1}{\sqrt{n}} Z + \frac{1}{\sqrt{n}} \Delta, \quad (66)$$

with Z, Δ defined as per the theorem statement. Further Z is Gaussian with the stated covariance because it is a linear function of the Gaussian vector $W \sim \mathbf{N}(0, \sigma^2 \mathbf{I}_{p \times p})$.

We are left with the task of proving the bound (11) on Δ . Note that by definition (2.1), we have

$$\|\Delta\|_\infty \leq \sqrt{n} |M \hat{\Sigma} - \mathbf{I}|_\infty \|\hat{\theta}^n - \theta_0\|_1 = \sqrt{n} \mu_* \|\hat{\theta}^n - \theta_0\|_1. \quad (67)$$

By [BvdG11, Theorem 6.1, Lemma 6.2], we have, for any $\lambda \geq 4\sigma\sqrt{2K \log(pe^{t^2/2})/n}$

$$\mathbb{P}\left(\|\hat{\theta}^n - \theta_0\|_1 \geq \frac{4\lambda s_0}{\phi_0^2}\right) \leq 2e^{-t^2/2}. \quad (68)$$

(More precisely, we consider the trivial generalization of [BvdG11, Lemma 6.2] to the case $(\mathbf{X}^T \mathbf{X}/n)_{ii} \leq K$, instead of $(\mathbf{X}^T \mathbf{X}/n)_{ii} = 1$ for all $i \in [p]$.)

Substituting Eq. (67) in the last bound, we get

$$\mathbb{P}\left(\|\Delta\|_\infty \geq \frac{4\lambda\mu_*s_0\sqrt{n}}{\phi_0^2}\right) \leq 2e^{-t^2/2}. \quad (69)$$

Finally, the claim follows by selecting t so that $e^{t^2/2} = p^{c_0}$.

6.2 Proof of Theorem 2.4.(a)

Note that the event \mathcal{E}_n requires two conditions. Hence, its complement

$$\mathcal{E}_n(\phi_0, s_0, K)^c = \mathcal{B}_{1,n}(\phi_0, s_0) \cup \mathcal{B}_{2,n}(K), \quad (70)$$

$$\mathcal{B}_{1,n}(\phi_0, s_0) \equiv \left\{ \mathbf{X} \in \mathbb{R}^{n \times p} : \min_{S: |S| \leq s_0} \phi(\widehat{\Sigma}, S) < \phi_0, \widehat{\Sigma} = (\mathbf{X}^T \mathbf{X}/n) \right\}, \quad (71)$$

$$\mathcal{B}_{2,n}(K) \equiv \left\{ \mathbf{X} \in \mathbb{R}^{n \times p} : \max_{i \in [p]} \widehat{\Sigma}_{i,i} \leq K, \widehat{\Sigma} = (\mathbf{X}^T \mathbf{X}/n) \right\}. \quad (72)$$

We will bound separately the probability of $\mathcal{B}_{1,n}$ and the probability of $\mathcal{B}_{2,n}$. The claim of Theorem 2.4.(a) follows by union bound.

6.2.1 Controlling $\mathcal{B}_{1,n}(\phi_0, s_0)$

It is also useful to recall the notion of restricted eigenvalue, introduced by Bickel, Ritov and Tsybakov [BRT09].

Definition 6.1. *Given a symmetric matrix $Q \in \mathbb{R}^{p \times p}$ an integer $s_0 \geq 1$, and $L > 0$, the restricted eigenvalue of Q is defined as*

$$\phi_{\text{RE}}^2(Q, s_0, L) \equiv \min_{S \subseteq [p], |S| \leq s_0} \min_{\theta \in \mathbb{R}^p} \left\{ \frac{\langle \theta, Q \theta \rangle}{\|\theta_S\|_2^2} : \theta \in \mathbb{R}^p, \|\theta_{S^c}\|_1 \leq L \|\theta_S\|_1 \right\}. \quad (73)$$

Rudelson and Zhou [RZ13] prove that, if the population covariance satisfies the restricted eigenvalue condition, then the sample covariance satisfies it as well, with high probability. More precisely [RZ13, Theorem 6], the following happens for some $c_* \leq 2000$, $m \equiv c_* s_0 C_{\max}^2 / \phi_{\text{RE}}^2(\Sigma, s_0, 9)$, and every $n \geq 4c_* m \kappa^4 \log(60ep/(m\kappa))$ we have

$$\mathbb{P}\left(\phi_{\text{RE}}(\widehat{\Sigma}, s_0, 3) \geq \frac{1}{2} \phi_{\text{RE}}(\Sigma, s_0, 9)\right) \geq 1 - 2e^{-n/(4c_* \kappa^4)} \quad (74)$$

Note that $\phi_{\text{RE}}(\Sigma, s_0, 9) \geq \sigma_{\min}(\Sigma)^{1/2} \geq C_{\min}^{1/2}$ and, by Cauchy-Schwartz $\min_{S: |S| \leq s_0} \phi(\widehat{\Sigma}, S) \geq \phi_{\text{RE}}(\widehat{\Sigma}, s_0, 3)$. With the definitions in the statement (cf. Eq. (13)), we therefore have

$$\mathbb{P}\left(\min_{S: |S| \leq s_0} \phi(\widehat{\Sigma}, S) \geq \frac{1}{2} C_{\min}^{1/2}\right) \geq 1 - 2e^{-c_1 n}. \quad (75)$$

Equivalently, $\mathbb{P}(\mathcal{B}_{1,n}(\phi_0, s_0)) \leq 2e^{-c_1 n}$.

6.2.2 Controlling $\mathcal{B}_{2,n}(K)$

By definition

$$\widehat{\Sigma}_{ii} - 1 = \frac{1}{n} \sum_{\ell=1}^n (\langle X_\ell, e_i \rangle^2 - 1) = \frac{1}{n} \sum_{\ell=1}^n u_\ell, \quad (76)$$

Note that u_ℓ are independent centered random variables. Further (recalling that, for any random variables U, V , $\|U + V\|_{\psi_1} \leq \|U\|_{\psi_1} + \|V\|_{\psi_1}$, and $\|U^2\|_{\psi_1} \leq 2\|U\|_{\psi_2}^2$) they are subexponential with subexponential norm

$$\|u_\ell\|_{\psi_1} \leq 2\|\langle X_\ell, e_i \rangle^2\|_{\psi_1} \leq 4\|\langle X_\ell, e_i \rangle\|_{\psi_1}^2 \quad (77)$$

$$\leq 4\|\langle \Sigma^{-1/2} X_\ell, \Sigma^{1/2} e_i \rangle\|_{\psi_1}^2 \quad (78)$$

$$\leq 4\kappa^2 \|\Sigma^{1/2} e_i\|_2^2 = 4\kappa^2 \Sigma_{ii} = 4\kappa^2. \quad (79)$$

By Bernstein-type inequality for centered subexponential random variables [Ver12], we get

$$\mathbb{P}\left\{\frac{1}{n} \left| \sum_{\ell=1}^n u_\ell \right| \geq \varepsilon\right\} \leq 2 \exp\left[-\frac{n}{6} \min\left(\left(\frac{\varepsilon}{4e\kappa^2}\right)^2, \frac{\varepsilon}{4e\kappa^2}\right)\right]. \quad (80)$$

Hence, for all ε such that $\varepsilon/(e\kappa^2) \in [\sqrt{(48 \log p)/n}, 4]$,

$$\mathbb{P}\left(\max_{i \in [p]} \widehat{\Sigma}_{ii} \geq 1 + \varepsilon\right) \leq 2p \exp\left(-\frac{n\varepsilon^2}{24e^2\kappa^4}\right) \leq 2e^{-c_1 n}, \quad (81)$$

which implies $\mathbb{P}(\mathbf{X} \in \mathcal{B}_{2,n}(K)) \leq 2e^{-c_1 n}$ for all $K - 1 \geq 20\kappa^2 \sqrt{(\log p)/n} \geq \sqrt{(48e^2\kappa^4 \log p)/n}$.

6.3 Proof of Theorem 2.4.(b)

Obviously, we have

$$\mu_{\min}(\mathbf{X}) \leq |\Sigma^{-1} \widehat{\Sigma} - \mathbf{I}|, \quad (82)$$

and hence the statement follows immediately from the following estimate.

Lemma 6.2. *Consider a random design matrix $\mathbf{X} \in \mathbb{R}^{p \times p}$, with i.i.d. rows having mean zero and population covariance Σ . Assume that*

(i) *We have $\sigma_{\min}(\Sigma) \geq C_{\min} > 0$, and $\sigma_{\max}(\Sigma) \leq C_{\max} < \infty$.*

(ii) *The rows of $X\Sigma^{-1/2}$ are sub-gaussian with $\kappa = \|\Sigma^{-1/2} X_1\|_{\psi_2}$.*

Let $\widehat{\Sigma} = (\mathbf{X}^\top \mathbf{X})/n$ be the empirical covariance. Then, for any constant $C > 0$, the following holds true.

$$\mathbb{P}\left\{\left|\Sigma^{-1} \widehat{\Sigma} - \mathbf{I}\right|_{\infty} \geq a \sqrt{\frac{\log p}{n}}\right\} \leq 2p^{-c_2}, \quad (83)$$

with $c_2 = (a^2 C_{\min}) / (24e^2 \kappa^4 C_{\max}) - 2$.

Proof of Lemma 6.2. The proof is based on Bernstein-type inequality for sub-exponential random variables [Ver12]. Let $\tilde{X}_\ell = \Sigma^{-1/2} X_\ell$, for $\ell \in [n]$, and write

$$Z \equiv \Sigma^{-1} \hat{\Sigma} - \mathbf{I} = \frac{1}{n} \sum_{\ell=1}^n \left\{ \Sigma^{-1} X_\ell X_\ell^\top - \mathbf{I} \right\} = \frac{1}{n} \sum_{\ell=1}^n \left\{ \Sigma^{-1/2} \tilde{X}_\ell \tilde{X}_\ell^\top \Sigma^{1/2} - \mathbf{I} \right\}.$$

Fix $i, j \in [p]$, and for $\ell \in [n]$, let $v_\ell^{(ij)} = \langle \Sigma_{i,\cdot}^{-1/2}, \tilde{X}_\ell \rangle \langle \Sigma_{j,\cdot}^{1/2}, \tilde{X}_\ell \rangle - \delta_{i,j}$, where $\delta_{i,j} = \mathbf{1}_{\{i=j\}}$. Notice that $\mathbb{E}(v_\ell^{(ij)}) = 0$, and the $v_\ell^{(ij)}$ are independent for $\ell \in [n]$. Also, $Z_{i,j} = (1/n) \sum_{\ell=1}^n v_\ell^{(ij)}$. By [Ver12, Remark 5.18], we have

$$\|v_\ell^{(ij)}\|_{\psi_1} \leq 2 \|\langle \Sigma_{i,\cdot}^{-1/2}, \tilde{X}_\ell \rangle \langle \Sigma_{j,\cdot}^{1/2}, \tilde{X}_\ell \rangle\|_{\psi_1}.$$

Moreover, for any two random variables X and Y , we have

$$\begin{aligned} \|XY\|_{\psi_1} &= \sup_{p \geq 1} p^{-1} \mathbb{E}(|XY|^p)^{1/p} \\ &\leq \sup_{p \geq 1} p^{-1} \mathbb{E}(|X|^{2p})^{1/2p} \mathbb{E}(|Y|^{2p})^{1/2p} \\ &\leq 2 \left(\sup_{q \geq 2} q^{-1/2} \mathbb{E}(|X|^q)^{1/q} \right) \left(\sup_{q \geq 2} q^{-1/2} \mathbb{E}(|Y|^q)^{1/q} \right) \\ &\leq 2 \|X\|_{\psi_2} \|Y\|_{\psi_2}. \end{aligned}$$

Hence, by assumption (ii), we obtain

$$\begin{aligned} \|v_\ell^{(ij)}\|_{\psi_1} &\leq 2 \|\langle \Sigma_{i,\cdot}^{-1/2}, \tilde{X}_\ell \rangle\|_{\psi_2} \|\langle \Sigma_{j,\cdot}^{1/2}, \tilde{X}_\ell \rangle\|_{\psi_2} \\ &\leq 2 \|\Sigma_{i,\cdot}^{-1/2}\|_2 \|\Sigma_{j,\cdot}^{1/2}\|_2 \kappa^2 \leq 2 \sqrt{C_{\max}/C_{\min}} \kappa^2. \end{aligned}$$

Let $\kappa' = 2\sqrt{C_{\max}/C_{\min}} \kappa^2$. Applying Bernstein-type inequality for centered sub-exponential random variables [Ver12], we get

$$\mathbb{P} \left\{ \frac{1}{n} \left| \sum_{\ell=1}^n v_\ell^{(ij)} \right| \geq \varepsilon \right\} \leq 2 \exp \left[- \frac{n}{6} \min \left(\left(\frac{\varepsilon}{e\kappa'} \right)^2, \frac{\varepsilon}{e\kappa'} \right) \right].$$

Choosing $\varepsilon = a\sqrt{(\log p)/n}$, and assuming $n \geq [a/(e\kappa')]^2 \log p$, we arrive at

$$\mathbb{P} \left\{ \frac{1}{n} \left| \sum_{\ell=1}^n v_\ell^{(ij)} \right| \geq a\sqrt{\frac{\log p}{n}} \right\} \leq 2p^{-a^2/(6e^2\kappa'^2)}.$$

The result follows by union bounding over all possible pairs $i, j \in [p]$. □

6.4 Proof of Theorem 2.5

Let

$$\Delta_0 \equiv \left(\frac{16ac\sigma}{C_{\min}} \right) \frac{s_0 \log p}{\sqrt{n}} \tag{84}$$

be a shorthand for the bound on $\|\Delta\|_\infty$ appearing in Eq. (17). Then we have

$$\begin{aligned} \mathbb{P}\left(\|\Delta\|_\infty \geq \Delta_0\right) &\leq \mathbb{P}\left(\{\|\Delta\|_\infty \geq \Delta_0\} \cap \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)\right) \\ &\quad + \mathbb{P}(\mathcal{E}_n^c(C_{\min}^{1/2}/2, s_0, 3/2)) + \mathbb{P}(\mathcal{G}_n^c(a)) \end{aligned} \quad (85)$$

$$\leq \mathbb{P}\left(\{\|\Delta\|_\infty \geq \Delta_0\} \cap \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)\right) + 4e^{-c_1 n} + 2p^{-c_2}, \quad (86)$$

where, in the first equation \mathcal{A}^c denotes the complement of event \mathcal{A} and the second inequality follows from Theorem 2.4. Notice, in particular, that the bound (13) can be applied for $K = 3/2$ since, under the present assumptions $20\kappa^2\sqrt{(\log p)/n} \leq 1/2$.

Finally

$$\begin{aligned} \mathbb{P}\left(\{\|\Delta\|_\infty \geq \Delta_0\} \cap \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)\right) \\ \leq \sup_{\mathbf{X} \in \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)} \mathbb{P}\left(\|\Delta\|_\infty \geq \Delta_0 \mid \mathbf{X}\right) \leq 2p^{-\tilde{c}_0}, \end{aligned} \quad (87)$$

Here the last inequality follows from Theorem 2.3 applied per given $\mathbf{X} \in \mathcal{E}_n(C_{\min}^{1/2}/2, s_0, 3/2) \cap \mathcal{G}_n(a)$ and hence using the bound (11) with $\phi_0 = C_{\min}^{1/2}/2$, $K = 3/2$, $\mu_* = a\sqrt{(\log p)/n}$.

6.5 Proof of Lemma 3.2

We will prove that, under the stated assumptions

$$\limsup_{n \rightarrow \infty} \sup_{\|\theta_0\|_0 \leq s_0} \mathbb{P}\left\{\frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \leq x\right\} \leq \Phi(x). \quad (88)$$

A matching lower bound follows by a completely analogous argument.

Notice that by Eq. (16), we have

$$\frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\sigma[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} = \frac{e_i^\top M\mathbf{X}^\top W}{\sigma[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} + \frac{\Delta_i}{\sigma[M\hat{\Sigma}M^\top]_{i,i}^{1/2}}. \quad (89)$$

Let $V = \mathbf{X}M^\top e_i / (\sigma[M\hat{\Sigma}M^\top]_{i,i}^{1/2})$ and $\tilde{Z} \equiv V^\top W$. We claim that $\tilde{Z} \sim \mathcal{N}(0, 1)$. To see this, note that $\|V\|_2 = 1$, and V and W are independent. Hence,

$$\mathbb{P}(\tilde{Z} \leq x) = \mathbb{E}\{\mathbb{P}(V^\top W \leq x \mid V)\} = \mathbb{E}\{\Phi(x) \mid V\} = \Phi(x), \quad (90)$$

which proves our claim. In order to prove Eq. (88), fix $\varepsilon > 0$ and write

$$\mathbb{P}\left(\frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \leq x\right) = \mathbb{P}\left(\frac{\sigma}{\hat{\sigma}}\tilde{Z} + \frac{\Delta_i}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq x\right) \quad (91)$$

$$\leq \mathbb{P}\left(\frac{\sigma}{\hat{\sigma}}\tilde{Z} \leq x + \varepsilon\right) + \mathbb{P}\left(\frac{|\Delta_i|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq \varepsilon\right) \quad (92)$$

$$\leq \mathbb{P}\left(\tilde{Z} \leq x + 2\varepsilon + \varepsilon|x\right) + \mathbb{P}\left(\frac{|\Delta_i|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq \varepsilon\right) + \mathbb{P}\left(\left|\frac{\hat{\sigma}}{\sigma} - 1\right| \geq \varepsilon\right). \quad (93)$$

By taking the limit and using the assumption (30), we obtain

$$\limsup_{n \rightarrow \infty} \sup_{\|\theta_0\|_0 \leq s_0} \mathbb{P} \left(\frac{\sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \leq x \right) \leq \tag{94}$$

$$\Phi(x + 2\varepsilon + \varepsilon|x|) + \limsup_{n \rightarrow \infty} \sup_{\|\theta_0\|_0 \leq s_0} \mathbb{P} \left(\frac{|\Delta_i|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq \varepsilon \right).$$

Since $\varepsilon > 0$ is arbitrary, it is therefore sufficient to show that the limit on the right hand side vanishes for any $\varepsilon > 0$.

Note that $[M\hat{\Sigma}M^\top]_{i,i} \geq 1/(4\hat{\Sigma}_{ii})$ for all n large enough, by Lemma 3.1, and since $\mu = a\sqrt{(\log p)/n} \rightarrow 0$ as $n, p \rightarrow \infty$. We have therefore

$$\mathbb{P} \left(\frac{|\Delta_i|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq \varepsilon \right) \leq \mathbb{P} \left(\frac{2}{\hat{\sigma}} \hat{\Sigma}_{ii}^{1/2} |\Delta_i| \geq \varepsilon \right) \tag{95}$$

$$\leq \mathbb{P} \left(\frac{8}{\sigma} |\Delta_i| \geq \varepsilon \right) + \mathbb{P} \left(\frac{\hat{\sigma}}{\sigma} \geq 2 \right) + \mathbb{P}(\hat{\Sigma}_{ii} \geq \sqrt{2}). \tag{96}$$

Note that $\mathbb{P}((\hat{\sigma}/\sigma) \geq 2) \rightarrow 0$ by assumption (30), and $\mathbb{P}(\hat{\Sigma}_{ii} \geq \sqrt{2}) \rightarrow 0$ by Theorem 2.4.(b). Hence

$$\limsup_{n \rightarrow \infty} \sup_{\|\theta_0\|_0 \leq s_0} \mathbb{P} \left(\frac{|\Delta_i|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \geq \varepsilon \right) \leq \limsup_{n \rightarrow \infty} \sup_{\|\theta_0\|_0 \leq s_0} \mathbb{P} \left(\|\Delta\|_\infty \geq \frac{\varepsilon\sigma}{8} \right) \tag{97}$$

$$\leq \limsup_{n \rightarrow \infty} (4e^{-c_1 n} + 4p^{-(\tilde{c}_0 \wedge c_2)}) = 0, \tag{98}$$

where the last inequality follows from Eq. (17) since $s_0 = o(\sqrt{n}/\log p)$ and hence $(16acs_0 \log p)/(C_{\min}\sqrt{n}) \leq \varepsilon/8$ for all n large enough.

This completes the proof of Eq. (88). The matching lower bound follows by the same argument.

6.6 Proof of Theorem 3.5

We begin with proving Eq. (42). Defining $Z_i \equiv \sqrt{n}(\hat{\theta}_i^u - \theta_{0,i})/(\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2})$, we have

$$\begin{aligned} \lim_{n \rightarrow \infty} \alpha_{i,n}(\hat{T}) &= \lim_{n \rightarrow \infty} \sup_{\theta_0} \left\{ \mathbb{P}(P_i \leq \alpha) : i \in [p], \|\theta_0\|_0 \leq s_0, \theta_{0,i} = 0 \right\} \\ &= \lim_{n \rightarrow \infty} \sup_{\theta_0} \left\{ \mathbb{P} \left(\Phi^{-1} \left(1 - \frac{\alpha}{2} \right) \leq \frac{\sqrt{n}|\hat{\theta}_i^u|}{\hat{\sigma}[M\hat{\Sigma}M^\top]_{i,i}^{1/2}} \right) : i \in [p], \|\theta_0\|_0 \leq s_0, \theta_{0,i} = 0 \right\} \\ &= \lim_{n \rightarrow \infty} \sup_{\theta_0} \left\{ \mathbb{P} \left(\Phi^{-1} \left(1 - \frac{\alpha}{2} \right) \leq |Z_i| \right) : i \in [p], \|\theta_0\|_0 \leq s_0 \right\} \leq \alpha, \end{aligned}$$

where the last inequality follows from Lemma 3.2.

We next prove Eq. (43). Recall that $\Sigma_{\cdot,i}^{-1}$ is a feasible solution of (4), for $1 \leq i \leq p$ with probability at least $1 - 2p^{-c_2}$, as per Lemma 6.2). On this event, letting m_i be the solution of the optimization

problem (4), we have

$$\begin{aligned}
m_i^\top \widehat{\Sigma} m_i &\leq \Sigma_{i,\cdot}^{-1} \widehat{\Sigma} \Sigma_{i,\cdot}^{-1} \\
&= (\Sigma_{i,\cdot}^{-1} \widehat{\Sigma} \Sigma_{i,\cdot}^{-1} - \Sigma_{ii}^{-1}) + \Sigma_{i,i}^{-1} \\
&= \frac{1}{n} \sum_{j=1}^N (V_j^2 - \Sigma_{ii}^{-1}) + \Sigma_{i,i}^{-1},
\end{aligned} \tag{99}$$

where $V_j = \Sigma_{i,\cdot}^{-1} X_j$ are i.i.d. with $\mathbb{E}(V_j^2) = \Sigma_{ii}^{-1}$ and sub-gaussian norm $\|V_j\|_{\psi_2} \leq \|\Sigma_{i,\cdot}^{-1/2}\|_2 \|\Sigma^{-1/2} X_j\|_{\psi_2} \leq \kappa \sqrt{\Sigma_{i,i}^{-1}}$. Letting $U_j = V_j^2 - \Sigma_{ii}^{-1}$, we have that U_j is zero mean and sub-exponential with $\|U_j\|_{\psi_1} \leq 2\|V_j^2\|_{\psi_1} \leq 2\|V_j\|_{\psi_2}^2 \leq 2\kappa^2 \Sigma_{ii}^{-1} \leq 2\kappa^2 \sigma_{\min}(\Sigma)^{-1} \leq 2\kappa^2 C_{\min}^{-1} \equiv \kappa'$. Hence, by applying Bernstein inequality (as, for instance, in the proof of Lemma 6.2), we have, for $\varepsilon \leq e\kappa'$,

$$\mathbb{P}\left(m_i^\top \widehat{\Sigma} m_i \geq \Sigma_{i,i}^{-1} + \varepsilon\right) \leq 2e^{-(n/6)(\varepsilon/e\kappa')^2} + 2p^{-c_2}. \tag{100}$$

Therefore, by Borel-Cantelli (since we can make $c_2 \geq 2$ by a suitable choice of a), we have, almost surely

$$\limsup_{n \rightarrow \infty} [m_i^\top \widehat{\Sigma} m_i - \Sigma_{i,i}^{-1}] \leq 0. \tag{101}$$

This bound leads to a lower bound for the power. First of all, a straightforward manipulation yields as follows, letting $z_* \equiv \Phi^{-1}(1 - \alpha/2)$:

$$\begin{aligned}
&\liminf_{n \rightarrow \infty} \frac{1 - \beta_{i,n}(\widehat{T}; \gamma)}{1 - \beta_{i,n}^*(\gamma)} \\
&= \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} \inf_{\theta_0} \left\{ \mathbb{P}(P_i \leq \alpha) : \|\theta_0\|_0 \leq s_0, |\theta_{0,i}| \geq \gamma \right\} \\
&= \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} \inf_{\theta_0} \left\{ \mathbb{P}\left(z_* \leq \frac{\sqrt{n}|\widehat{\theta}_i^u|}{\widehat{\sigma}[M\widehat{\Sigma}M^\top]_{i,i}^{1/2}}\right) : \|\theta_0\|_0 \leq s_0, |\theta_{0,i}| \geq \gamma \right\} \\
&= \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} \inf_{\theta_0} \left\{ \mathbb{P}\left(z_* \leq \left|Z_i + \frac{\sqrt{n}\theta_{0,i}}{\widehat{\sigma}[M\widehat{\Sigma}M^\top]_{i,i}^{1/2}}\right|\right) : \|\theta_0\|_0 \leq s_0, |\theta_{0,i}| \geq \gamma \right\} \\
&\stackrel{(a)}{\geq} \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} \inf_{\theta_0} \left\{ \mathbb{P}\left(z_* \leq \left|Z_i + \frac{\sqrt{n}\gamma}{\sigma[\Sigma_{i,i}^{-1}]^{1/2}}\right|\right) : \|\theta_0\|_0 \leq s_0 \right\} \\
&= \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} \left\{ 1 - \Phi\left(z_* - \frac{\sqrt{n}\gamma}{\sigma[\Sigma_{i,i}^{-1}]^{1/2}}\right) + \Phi\left(-z_* - \frac{\sqrt{n}\gamma}{\sigma[\Sigma_{i,i}^{-1}]^{1/2}}\right) \right\} \\
&= \liminf_{n \rightarrow \infty} \frac{1}{1 - \beta_{i,n}^*(\gamma)} G\left(\alpha, \frac{\sqrt{n}\gamma}{\sigma[\Sigma_{i,i}^{-1}]^{1/2}}\right) = 1.
\end{aligned}$$

Here (a) follows from Eq. (101) and the fact $|\theta_{0,i}| \geq \gamma$.

6.7 Proof of Theorem 4.1

Under the assumptions of Theorem 2.5 and assuming $s_0 = o(\sqrt{n}/\log p)$, we have

$$\sqrt{n}(\widehat{\theta}^u - \theta_0) = \frac{1}{\sqrt{n}}M\mathbf{X}^\top W + \Delta$$

with $\|\Delta\|_\infty = o(1)$. Using Lemma 3.1, we have

$$\frac{\sqrt{n}(\widehat{\theta}_i^u - \theta_{0,i})}{\sigma[M\widehat{\Sigma}M^\top]_{i,i}^{1/2}} = Z_i + o(1), \quad \text{with } Z_i \equiv \frac{1}{\sqrt{n}} \frac{m_i^\top \mathbf{X}^\top W}{\sigma[m_i^\top \widehat{\Sigma} m_i]^{1/2}}.$$

The following lemma characterizes the limiting distribution of $Z_i|\mathbf{X}$ which implies the validity of the proposed p -value P_i and confidence intervals.

Lemma 6.3. *Suppose that the noise variables W_i are independent with $\mathbb{E}(W_i) = 0$, and $\mathbb{E}(W_i^2) = \sigma^2$, and $E(|W_i|^{2+a}) \leq C\sigma^{2+a}$ for some $a > (1/2 - \beta)^{-1}$. Let $M = (m_1, \dots, m_p)^\top$ be the matrix with rows m_i^\top obtained by solving optimization problem (57). For $i \in [p]$, define*

$$Z_i = \frac{1}{\sqrt{n}} \frac{m_i^\top \mathbf{X}^\top W}{\sigma[m_i^\top \widehat{\Sigma} m_i]^{1/2}}.$$

Under the assumptions of Theorem 2.5, for any sequence $i = i(n) \in [p]$, and any $x \in \mathbb{R}$, we have

$$\lim_{n \rightarrow \infty} \mathbb{P}(Z_i \leq x | \mathbf{X}) = \Phi(x).$$

Lemma 6.3 is proved in Appendix A.2.

Acknowledgments

A.J. is supported by a Caroline and Fabian Pease Stanford Graduate Fellowship. This work was partially supported by the NSF CAREER award CCF-0743978, the NSF grant DMS-0806211, and the grants AFOSR/DARPA FA9550-12-1-0411 and FA9550-13-1-0036.

A Proof of technical lemmas

A.1 Proof of Lemma 3.1

Let $C_i(\mu)$ be the optimal value of the optimization problem (4). We claim that

$$C_i(\mu) \geq \frac{(1-\mu)^2}{\widehat{\Sigma}_{ii}}. \quad (102)$$

To prove this claim notice that the constraint implies (by considering its i -th component):

$$1 - \langle e_i, \widehat{\Sigma}m \rangle \leq \mu.$$

Therefore if \tilde{m} is feasible and $c \geq 0$, then

$$\langle \tilde{m}, \widehat{\Sigma}\tilde{m} \rangle \geq \langle \tilde{m}, \widehat{\Sigma}\tilde{m} \rangle + c(1-\mu) - c\langle e_i, \widehat{\Sigma}\tilde{m} \rangle \geq \min_m \left\{ \langle m, \widehat{\Sigma}m \rangle + c(1-\mu) - c\langle e_i, \widehat{\Sigma}m \rangle \right\}.$$

Minimizing over all feasible \tilde{m} gives

$$C_i(\mu) \geq \min_m \left\{ \langle m, \widehat{\Sigma}m \rangle + c(1-\mu) - c\langle e_i, \widehat{\Sigma}m \rangle \right\}. \quad (103)$$

The minimum over m is achieved at $m = ce_i/2$. Plugging in for m , we get

$$C_i(\mu) \geq c(1-\mu) - \frac{c^2}{4}\widehat{\Sigma}_{ii} \quad (104)$$

Optimizing this bound over c , we obtain the claim (102), with the optimal choice being $c = 2(1-\mu)/\widehat{\Sigma}_{ii}$.

A.2 Proof of Lemma 6.3

Write

$$Z_i = \frac{1}{\sqrt{n}} \sum_{j=1}^n \xi_j \quad \text{with} \quad \xi_j \equiv \frac{m_i^\top X_j W_j}{\sigma[m_i^\top \widehat{\Sigma}m_i]^{1/2}}.$$

Conditional on \mathbf{X} , the summands ξ_j are zero mean and independent. Furthermore, $\sum_{j=1}^n \mathbb{E}(\xi_j^2 | \mathbf{X}) = n$. We next prove the Lindenberg condition as per Eq. (56). Let $c_n \equiv (m_i^\top \widehat{\Sigma}m_i)^{1/2}$. By Lemma 3.1, we have, almost surely, $\liminf_{n \rightarrow \infty} c_n \geq c_\infty > 0$. If all the optimization problems in (57) are feasible, then $|\xi_j| \leq c_n^{-1} \|\mathbf{X}m_i\|_\infty \|W\|_\infty / \sigma \leq c_n^{-1} n^\beta (\|W\|_\infty / \sigma)$. Hence,

$$\begin{aligned} \lim_{n \rightarrow \infty} \frac{1}{n} \sum_{j=1}^n \mathbb{E} \left(\xi_j^2 \mathbb{I}_{\{|\xi_j| > \varepsilon \sqrt{n}\}} | \mathbf{X} \right) &\leq \lim_{n \rightarrow \infty} \frac{1}{n} \sum_{j=1}^n \mathbb{E} \left(\xi_j^2 \mathbb{I}_{\{\|W\|_\infty / \sigma > \varepsilon c_n n^{1/2-\beta}\}} | \mathbf{X} \right) \\ &= \lim_{n \rightarrow \infty} \frac{1}{n} \sum_{j=1}^n \frac{m_i^\top X_j X_j^\top m_i}{m_i^\top \widehat{\Sigma}m_i} \mathbb{E}(\widetilde{W}_j^2 \mathbb{I}_{\{\|\widetilde{W}\|_\infty > \varepsilon c_\infty n^{1/2-\beta}\}}) \\ &\leq \lim_{n \rightarrow \infty} n \mathbb{E}(\widetilde{W}_1^2 \mathbb{I}_{\{|\widetilde{W}_1| > \varepsilon c_\infty n^{1/2-\beta}\}}) \\ &\leq c'(\varepsilon) \lim_{n \rightarrow \infty} n^{1-a(1/2-\beta)} \mathbb{E}\{|\widetilde{W}_1|^{2+a}\} = 0. \end{aligned}$$

where $\widetilde{W}_j = W_j/\sigma$ and the last limit follows by taking $a > (1/2 - \beta)^{-1}$ as per the assumptions.

Using Lindenberg central limit theorem, we obtain $Z_i|\mathbf{X}$ converges weakly to standard normal distribution, and hence, \mathbf{X} -almost surely

$$\lim_{n \rightarrow \infty} \mathbb{P}(Z_i \leq x | \mathbf{X}) = \Phi(x).$$

What remains is to show that with high probability all the p optimization problems in (57) are feasible. In particular, we show that $\Sigma_{i,\cdot}^{-1}$ is a feasible solution to the i -th optimization problem, for $i \in [p]$. By Lemma 6.2, $|\Sigma^{-1}\widehat{\Sigma} - \mathbf{I}|_\infty \leq \mu$, with high probability. Moreover,

$$\begin{aligned} \sup_{j \in [p]} \|\Sigma_{i,\cdot}^{-1} X_j\|_{\psi_2} &= \sup_{j \in [p]} \|\Sigma_{i,\cdot}^{-1/2} \Sigma^{-1/2} X_j\|_{\psi_2} \\ &= \|\Sigma_{i,\cdot}^{-1/2}\|_2 \sup_{j \in [p]} \|\Sigma^{-1/2} X_j\|_{\psi_2} \\ &= [\Sigma_{i,i}^{-1}]^{1/2} \sup_{j \in [p]} \|\Sigma^{-1/2} X_j\|_{\psi_2} = O(1). \end{aligned}$$

Using tail bound for sub-gaussian variables $\Sigma_{i,\cdot}^{-1} X_j$ and union bounding over $j \in [n]$, we get

$$\mathbb{P}(\|\mathbf{X}\Sigma_{i,\cdot}^{-1}\|_\infty > n^\beta) \leq ne^{-cn^{2\beta}},$$

for some constant $c > 0$. Note that $s_0 = o(\sqrt{n}/\log p)$ implies $p = e^{o(n^{2\beta})}$. Hence, eventually almost surely, $\Sigma_{i,\cdot}^{-1}$ is a feasible solution to optimization problem (57), for all $i \in [p]$.

B Corollaries of Theorem 2.5

B.1 Proof of Corollary 2.7

By Theorem 2.3, for any $\mathbf{X} \in \mathcal{E}_n(\sqrt{C_{\min}}/2, s_0, 3/2) \cap \mathcal{G}_n(a)$, we have

$$\mathbb{P}\left\{\|\Delta\|_\infty \geq Lc \mid \mathbf{X}\right\} \leq 2p^{1-(c^2/48)}, \quad L \equiv \frac{16a\sigma}{C_{\min}} \frac{s_0 \log p}{\sqrt{n}}. \quad (105)$$

(This is obtained by setting $\phi_0 = C_{\min}^{1/2}/2$, $K = 3/2$, $\mu_* = a\sqrt{(\log p)/n}$ in Eq. (11). Hence

$$\|\text{Bias}(\widehat{\theta}^u)\|_\infty \leq \frac{1}{\sqrt{n}} \mathbb{E}\{\|\Delta\|_\infty \mid \mathbf{X}\} \quad (106)$$

$$= \frac{L'}{\sqrt{n}} \int_0^\infty \mathbb{P}\left\{\|\Delta\|_\infty \geq B'c \mid \mathbf{X}\right\} dc \quad (107)$$

$$\leq \frac{L}{\sqrt{n}} \int_0^\infty \min(1, p^{1-(c^2/48)}) dc \leq \frac{10L}{\sqrt{n}}, \quad (108)$$

which coincides with Eq. (21). The probability estimate (22) simply follows from Theorem 2.4 using union bound.

B.2 Proof of Corollary 2.8

By Theorem 2.4.(a), we have (setting $C_{\min} = C_{\max} = \kappa = 1$):

$$\mathbb{P}\left(\mathbf{X} \in \mathcal{E}_n(1/2, s_0, 3/2)\right) \geq 1 - 4e^{-n/c_*}. \quad (109)$$

Further, by Lemma 6.2, with $\widehat{\Sigma} \equiv \mathbf{X}^\top \mathbf{X}/n$, we have

$$\mathbb{P}\left(\mu_*(\mathbf{X}; \mathbf{I}) \leq 30\sqrt{\frac{\log p}{n}}\right) \geq 1 - 2p^{-3}. \quad (110)$$

Finally, by an obvious consequence of the proof of Theorem 2.4.(a)

$$\mathbb{P}\left(\left\{\mathbf{X} : \min_{i \in [p]} \widehat{\Sigma}_{ii} \geq \frac{1}{2}\right\}\right) \geq 1 - 2e^{-n/c_*}. \quad (111)$$

Hence, defining

$$\mathcal{B}_n \equiv \mathcal{E}_n(1/2, s_0, 3/2) \cap \left\{\mathbf{X} \in \mathbb{R}^{n \times p} : \mu_*(\mathbf{X}; \mathbf{I}) \leq 30\sqrt{\frac{\log p}{n}}\right\} \cap \left\{\mathbf{X} : \min_{i \in [p]} \widehat{\Sigma}_{ii} \geq \frac{1}{2}\right\} \quad (112)$$

we have the desired probability bound (24).

Let

$$\widehat{\theta}^* \equiv \widehat{\theta}^n + \frac{1}{n} \mathbf{X}^\top (Y - \mathbf{X} \widehat{\theta}^n), \quad (113)$$

where $\widehat{\theta}^n = \widehat{\theta}^n(Y, \mathbf{X}; \lambda)$ is the LASSO solution with $\lambda = \sigma\sqrt{(c^2 \log p)/n}$. By Theorem 2.3, we have, for any $\mathbf{X} \in \mathcal{B}_n$

$$\widehat{\theta}^* = \theta_0 + \frac{1}{\sqrt{n}} Z + \frac{1}{\sqrt{n}} \Delta, \quad Z | \mathbf{X} \sim \mathbf{N}(0, \sigma^2 \widehat{\Sigma}), \quad (114)$$

and further

$$\mathbb{P}\left\{\|\Delta\|_\infty \geq \frac{240c\sigma s_0 \log p}{\sqrt{n}} \mid \mathbf{X}\right\} \leq 2p^{1-(c^2/48)}, \quad (115)$$

whence, proceeding as in the proof in the last section, we get, for some universal numerical constant c_{**} ,

$$\|\text{Bias}(\widehat{\theta}^u)\|_\infty \leq \frac{1}{\sqrt{n}} \mathbb{E}\left\{\|\Delta\|_\infty \mid \mathbf{X}\right\} \leq c_{**} \sigma \frac{s_0 \log p}{n}. \quad (116)$$

Next by Eq. (29) we have

$$\|\text{Bias}(\widehat{\theta}^n)\|_\infty \geq \left| \lambda \|\mathbb{E}\{v(\widehat{\theta}^n) | \mathbf{X}\}\|_\infty - \|\text{Bias}(\widehat{\theta}^u)\|_\infty \right|. \quad (117)$$

Hence, in order to prove Eq. (23), it is sufficient to prove that $\|\mathbb{E}\{v(\widehat{\theta}^n) | \mathbf{X}\}\|_\infty \geq 2/3$.

Note that $v(\widehat{\theta}^n)_i = 1$ whenever $\widehat{\theta}_i^n > 0$ and, and $|v(\widehat{\theta}^n)_i| \leq 1$, and therefore (letting $b_0 \equiv c_{**}\sigma(s_0 \log p)/n$)

$$1 - \mathbb{E}\{v(\widehat{\theta}^n)_i | \mathbf{X}\} \leq 2\mathbb{P}\left(\widehat{\theta}_i^n \leq 0 \mid \mathbf{X}\right) \leq 2\mathbb{P}\left(\widehat{\theta}_i^n \leq \lambda \mid \mathbf{X}\right) \quad (118)$$

$$\leq 2\mathbb{P}\left(\theta_{0,i} + \frac{1}{\sqrt{n}}Z_i + \frac{1}{\sqrt{n}}\Delta_i \leq \lambda \mid \mathbf{X}\right) \quad (119)$$

$$\leq 2\mathbb{P}\left(\frac{1}{\sqrt{n}}Z_i \leq \lambda + b_0 - \theta_{0,i} \mid \mathbf{X}\right) \quad (120)$$

$$= 2\Phi\left((\lambda + b_0 - \theta_{0,i})\sqrt{n/(\sigma^2\widehat{\Sigma}_{ii})}\right) \quad (121)$$

$$\leq 2\Phi\left((\lambda + b_0 - \theta_{0,i})\sqrt{2n/(3\sigma^2)}\right) \quad (122)$$

with $\Phi(x)$ the standard normal distribution function, and in the last inequality we used the fact that $\max_{i \in [p]} \widehat{\Sigma}_{ii} \leq 3/2$ on \mathcal{B}_n . We then choose θ_0 so that $\theta_{0,i} \geq b_0 + \lambda + \sqrt{30\sigma^2/n}$, for $i \in [p]$ in the support of θ_0 . We therefore obtain

$$\mathbb{E}\{v(\widehat{\theta}^n)_i | \mathbf{X}\} \geq 1 - 2\Phi(-\sqrt{20}) \geq \frac{2}{3}. \quad (123)$$

This finishes the proof of Eq. (23). Equations (26) and (27) are obtained by substituting $\lambda = c\sigma\sqrt{(\log p)/n}$ and using Eq. (23).

C Proof of Lemma 3.3

Let $\mathcal{E}_n = \mathcal{E}_n(\phi_0, s_0, K)$ be the event defined as per Theorem 2.4.(a). In particular, we take $\phi_0 = C_{\min}^{1/2}/2$, and $K \geq 1 + 20\kappa^2\sqrt{(\log p)/n}$ (for, instance $K = 1.1$ will work for all n large enough since $(s_0 \log p)^2/n \rightarrow 0$, with $s_0 \geq 1$, by assumption). Further note that we can assume without loss of generality $n \geq \nu_0 s_0 \log(p/s_0)$, since $s_0 = o(\sqrt{n}/\log p)$. Fixing $\varepsilon > 0$, we have therefore

$$\mathbb{P}\left(\left|\frac{\widehat{\sigma}}{\sigma} - 1\right| \geq \varepsilon\right) \leq \sup_{\mathbf{X} \in \mathcal{E}_n} \mathbb{P}\left(\left|\frac{\widehat{\sigma}}{\sigma} - 1\right| \geq \varepsilon \mid \mathbf{X}\right) + \mathbb{P}(\mathbf{X} \notin \mathcal{E}_n) \quad (124)$$

$$\leq \sup_{\mathbf{X} \in \mathcal{E}_n} \mathbb{P}\left(\left|\frac{\widehat{\sigma}}{\sigma} - 1\right| \geq \varepsilon \mid \mathbf{X}\right) + 4e^{-c_1 n}, \quad (125)$$

where $c_1 > 0$ is a constant defined as per Theorem 2.4.(a).

We are therefore left with the task of bounding the first term in the last expression above, uniformly over $\theta_0 \in \mathbb{R}^p$, $\|\theta_0\|_0 \leq s_0$. For $\mathbf{X} \in \mathcal{E}_n$, we can apply [SZ12, Theorem 1.2] whereby (using the notations of [SZ12], with their λ_0 replaced by $\widetilde{\lambda}$) $\xi = 3$, $T = \text{supp}(\theta_0)$, $\kappa(\xi, T) \geq \phi_0$, $\eta_*(\widetilde{\lambda}, \xi) \leq 4s_0\widetilde{\lambda}^2/\phi_0^2$. By a straightforward manipulation of Eq. (13) in [SZ12], we have, for $\mathbf{X} \in \mathcal{E}$, and $\|\mathbf{X}^\top W/n\|_\infty \leq \widetilde{\lambda}/4$ (letting σ^* the oracle estimator of σ introduced there)

$$\left|\frac{\widehat{\sigma}}{\sigma^*} - 1\right| \leq \frac{4\sqrt{s_0}\widetilde{\lambda}}{\phi_0} \leq \frac{\varepsilon}{2} \quad (126)$$

where the last inequality follows for all n large enough since $s_0 = o(\sqrt{n}/\log p)$.

Hence

$$\sup_{\mathbf{X} \in \mathcal{E}_n} \mathbb{P} \left(\left| \frac{\widehat{\sigma}}{\sigma} - 1 \right| \geq \varepsilon \mid \mathbf{X} \right) \leq \sup_{\mathbf{X} \in \mathcal{E}_n} \mathbb{P} \left(\|\mathbf{X}^\top W/n\|_\infty \leq \tilde{\lambda}/4 \mid \mathbf{X} \right) + \sup_{\mathbf{X} \in \mathcal{E}_n} \mathbb{P} \left(\left| \frac{\sigma^*}{\sigma} - 1 \right| \geq \frac{1}{10} \mid \mathbf{X} \right), \quad (127)$$

where we note that the right hand side is independent of θ_0 . The first term vanishes as $n \rightarrow \infty$ by a standard tail bound on the supremum of p Gaussian random variables. The second term also vanishes because it is controlled by the tail of a chi-squared random variable [SZ12].

D Proof of Theorem 3.9

By definition, letting $\mathcal{F}_{p,s_0} \equiv \{x \in \mathbb{R}^p : \|x\|_0 \leq s_0\}$, and fixing $\varepsilon \in (0, 1/10)$

$$\text{FWER}(\widehat{T}^F, n) = \sup_{\theta_0 \in \mathcal{F}_{p,s_0}} \mathbb{P} \left\{ \exists i \in [p] \setminus \text{supp}(\theta_0) \text{ s.t. } \frac{\sqrt{n} |\widehat{\theta}_i^u - \theta_{0,i}|}{\widehat{\sigma} [M \widehat{\Sigma} M^\top]_{i,i}^{1/2}} \geq \Phi^{-1} \left(1 - \frac{\alpha}{2p} \right) \right\} \quad (128)$$

$$\begin{aligned} &\leq \sup_{\theta_0 \in \mathcal{F}_{p,s_0}} \mathbb{P} \left\{ \exists i \in [p] \setminus \text{supp}(\theta_0) \text{ s.t. } \frac{\sqrt{n} |\widehat{\theta}_i^u - \theta_{0,i}|}{\sigma [M \widehat{\Sigma} M^\top]_{i,i}^{1/2}} \geq (1 - \varepsilon) \Phi^{-1} \left(1 - \frac{\alpha}{2p} \right) \right\} \quad (129) \\ &\quad + \sup_{\theta_0 \in \mathcal{F}_{p,s_0}} \mathbb{P} \left(\left| \frac{\widehat{\sigma}}{\sigma} - 1 \right| \geq \frac{\varepsilon}{2} \right). \end{aligned}$$

Since the second term vanishes as $n \rightarrow \infty$ by assumption Eq. (30), it is sufficient to consider the first term. Using Bonferroni inequality, letting $z_\alpha(\varepsilon) \equiv (1 - \varepsilon) \Phi^{-1} \left(1 - \frac{\alpha}{2p} \right)$, we have

$$\limsup_{n \rightarrow \infty} \text{FWER}(\widehat{T}^F, n) \leq \limsup_{n \rightarrow \infty} \sum_{i=1}^p \sup_{\theta_0 \in \mathcal{F}_{p,s_0}, \theta_{0,i}=0} \mathbb{P} \left\{ \frac{\sqrt{n} |\widehat{\theta}_i^u - \theta_{0,i}|}{\sigma [M \widehat{\Sigma} M^\top]_{i,i}^{1/2}} \geq z_\alpha(\varepsilon) \right\} \quad (130)$$

$$= \limsup_{n \rightarrow \infty} \sum_{i=1}^p \sup_{\theta_0 \in \mathcal{F}_{p,s_0}, \theta_{0,i}=0} \mathbb{P} \left\{ \left| \widetilde{Z}_i + \frac{\Delta_i}{\sigma [M \widehat{\Sigma} M^\top]_{i,i}^{1/2}} \right| \geq z_\alpha(\varepsilon) \right\} \quad (131)$$

where, by Theorem 2.5, $\widetilde{Z}_i \sim \mathcal{N}(0, 1)$ and Δ_i is given by Eq. (16). We then have

$$\begin{aligned} \limsup_{n \rightarrow \infty} \text{FWER}(\widehat{T}^F, n) &\leq \limsup_{n \rightarrow \infty} \sum_{i=1}^p \mathbb{P} \{ |\widetilde{Z}_i| \geq z_\alpha(\varepsilon) - \varepsilon \} \\ &\quad + \limsup_{n \rightarrow \infty} \sum_{i=1}^p \sup_{\theta_0 \in \mathcal{F}_{p,s_0}, \theta_{0,i}=0} \mathbb{P} \left\{ \|\Delta\|_\infty \geq \frac{\varepsilon \sigma}{2 \widehat{\Sigma}_{ii}^{1/2}} \right\} \\ &\leq 2(1 - \Phi(z_\alpha(\varepsilon) - \varepsilon)) + \limsup_{n \rightarrow \infty} p \max_{i \in [p]} \mathbb{P}(\widehat{\Sigma}_{ii} \geq 2) \\ &\quad + \limsup_{n \rightarrow \infty} \sup_{\theta_0 \in \mathcal{F}_{p,s_0}, \theta_{0,i}=0} p \mathbb{P} \left\{ \|\Delta\|_\infty \geq \frac{\varepsilon \sigma}{4} \right\}, \quad (132) \end{aligned}$$

where in the first inequality, we used $[M \widehat{\Sigma} M^\top]_{i,i} \geq 1/(4 \widehat{\Sigma}_{ii})$ for all n large enough, by Lemma 3.1, and since $\mu = a \sqrt{(\log p)/n} \rightarrow 0$ as $n, p \rightarrow \infty$. Now the second term in the right hand side of Eq. (132)

vanishes by Theorem 2.4.(a), and the last term is zero by Theorem 2.5 since $\sqrt{n}/\log(p) \geq s_0 \geq 1$. Therefore

$$\limsup_{n \rightarrow \infty} \text{FWER}(\widehat{T}^F, n) \leq 2(1 - \Phi(z_\alpha(\varepsilon) - \varepsilon)), \quad (133)$$

and the claim follows by letting $\varepsilon \rightarrow 0$.

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