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Abstract

In high-dimensional regression scenarios with low signal-to-noise ratios, we assess the predictive performance of several machine learning algorithms. Theoretical insights show Ridge regression’s superiority in exploiting weak signals, surpassing a zero benchmark. In contrast, Lasso fails to exceed this baseline, indicating its learning limitations. Simulations reveal that Random Forest generally outperforms Gradient Boosted Regression Trees when signals are weak. Moreover, Neural Networks with $\ell_2$-regularization excel in capturing nonlinear functions of weak signals. Our empirical analysis across six economic datasets suggests that the weakness of signals, not necessarily the absence of sparsity, may be Lasso’s major limitation in economic predictions.

Keywords: Weak Signals, Precise Error, Machine Learning, Bayes Risk

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1 Introduction

In regression analysis, covariates with non-zero coefficients are recognized as true signals, while those with zero coefficients are considered false signals. In a population model, this distinction is clear-cut, resembling a “black and white” scenario. However, in finite samples, the presence of minuscule non-zero coefficients introduces a “gray” area, blurring the lines between true and false signals. This gray area represents weak signals—covariates that exert negligible influence on the outcome variable.

The investigation of weak signals holds tangible implications for economic and financial decision-making. Often, it is the collective impact of these weak signals that drives the outcomes in these fields. Supporting this, Figure 1 offers an empirical perspective, showcasing $R^2$ values gathered from a compendium of Economics and Finance journal articles published in 2022. The 25% quantiles of these $R^2$ values stand at 9.7% for economics and 5.8% for finance, suggesting that models in these disciplines frequently rely on covariates with modest explanatory power. Moreover, Figure 1 focuses solely on published papers, which likely skews towards studies with higher $R^2$ values due to selection bias. This suggests that the presence of weak signals may be even more widespread than the data here indicates.

The decision to incorporate weak signals into a regression model is fraught with the peril of overfitting, which can undermine predictive performance. This issue arises when the errors associated with estimating the coefficients of these weak signals outweigh the benefits of reducing bias that their inclusion offers. To include these variables or not thereby hinges on a trade-off between bias and variance. Compounding this challenge is the increasing prevalence of high-dimensional covariates in data-rich environments, a scenario frequently encountered recently, which can further exacerbate prediction errors due to the scarcity in terms of the sample size relative to the dimensionality of covariates.

Machine learning methods, known for their emphasis on variable selection and dimension reduction, have proven effective in mitigating overfitting and detecting true signals from false ones, particularly when the true signals are strong. These methods employ regularization techniques, such as penalizing the $\ell_1$ or $\ell_2$ norms of model parameters, to achieve this objective. A pivotal question arises: Can machines learn weak signals, or in other words, can they surpass the naive zero-estimator? The zero estimator, designed to ignore all covariates, serves as a passive baseline in the context of weak signals. If an estimator manages to surpass this baseline, it implies that it has effectively learned valuable signals. Conversely, failing to

\footnote{The comparison of the magnitudes of regression coefficients becomes meaningful only when the predictor variables have been normalized. This premise is implicitly assumed in our subsequent discussion.}
do so indicates a deficiency in its learning capacity.

In view of these considerations, we shift our focus to evaluate the relative performance of regularization techniques, particularly Ridge and Lasso estimators, against the zero estimator within high-dimensional regression contexts, where the dependent variable is driven by predictors that exhibit weak correlations with it.

In scenarios with sufficiently strong signals, both Lasso and Ridge estimators are expected to outperform the zero estimator by effectively capturing and utilizing at least some of these signals. Hence, accurately defining the notion of “weak” signals is crucial at the outset of our investigation. This definition serves a dual purpose: it prevents the scenario from defaulting to trivial comparisons akin to strong signal cases and ensures practical relevance to finite sample scenarios. We characterize a weak signal scenario as one in which the zero estimator, which always predicts the value zero, achieves the minimal Bayes prediction risk asymptotically. This setting turns out to encompass a sufficiently wide class of data generating processes (DGPs), and it approximates a finite sample reality in which the zero predictor serves as a competitive benchmark. If the zero predictor were not optimal, there would exist an estimator that dominates it, having identified some of the useful signals. In such instances, we would classify these scenarios as cases of strong signals.
In the defined weak signal scenarios, conventional error-bound analyses are insufficient in distinguishing the performance of different estimators. All three estimators — Lasso, Ridge, and the zero estimator — can attain the optimal Bayes prediction error, rendering them asymptotically indistinguishable. Intuitively, this occurs as the tuning parameters for both Ridge and Lasso estimators approach infinity, where Ridge tends toward zero, while Lasso becomes equivalent to zero.

To compare the predictive performance of these estimators, we have employed a precise error analysis approach, which enables us to zoom in and explicitly characterize the relative differences in the asymptotic behavior of zero, Ridge, and Lasso. This nuanced analysis reveals that with an appropriately chosen tuning parameter, Ridge asymptotically outperforms the zero estimator. In contrast, Lasso does not surpass the zero estimator’s performance, regardless of its tuning parameter choice. This finding aligns with the intuition that shrinkage methods like Ridge are more effective in environments with more homogenous signal strength. On the other hand, selection methods like Lasso are preferable in scenarios where there is a clear distinction between true and false signals. In weak signal contexts where this distinction is blurred, the advantage of Lasso tends to wane.

The assumptions about the DGPs in our analysis are sufficiently versatile to encompass a range of scenarios, including spike-and-slab models (George and McCulloch (1993)), where traditionally, Lasso has been the method of choice. Moreover, our analysis goes beyond the cases where regression coefficients follow a Gaussian distribution, a scenario where Ridge regression represents the posterior mean from a Bayesian perspective. This generality in DGP ensures that our findings do not inherently favor one estimator over the other, providing a balanced evaluation of their respective capabilities and limitations in weak signal scenarios.

Our study further emphasizes the validity of the cross-validation algorithm in identifying the optimal tuning parameter for Ridge regression, even in contexts with weak signals. This suggests that cross-validation remains a robust tool for model tuning, resilient to variations in signal strength. Moreover, we find that in the optimal Ridge regression, the out-of-sample $R^2$, a metric frequently used to evaluate the performance of different estimators on unseen data, proves a relevant indicator of the signal-to-noise ratio in the DGP, despite a notable gap between its asymptotic limit and the population $R^2$ inherent to the underlying regression.

In the final aspect of our theoretical analysis, we expand our framework to include models featuring a mix of signal strengths. This section specifically addresses scenarios in which a benchmark model contains potentially strong signals. We then shift our focus to assess benefit in harnessing predictive power from the weak signals that remain. To this end, we
derive ordinary least squares (OLS) residuals, from which the impact of the covariates in the benchmark model has been removed. Consistent with our earlier findings, applying Ridge regression to these residuals, using the remaining covariates, enhances predictive performance compared to a baseline estimator that ignores these additional covariates.

Our simulation analysis corroborates our theoretical findings: the Ridge estimator surpasses zero, which in turn edges out Lasso, especially in DGPs characterized by low $R^2$ values. Moving to more sophisticated machine learning techniques, we find that Random Forest (RF), yielding a dense model with almost all variables included, outperforms the zero estimator, which itself surpasses Gradient Boosted Regression Trees (GBRT). Resembling Lasso, GBRT tends to produce more sparse models in these scenarios of weak signal strength. Furthermore, Neural Networks (NNs), when integrated with the $\ell_2$-norm regularization, can effectively learn an unknown function of weak signals and yield superior predictions. In contrast, the application of an $\ell_1$-penalty in these networks does not yield comparable results.

From an empirical standpoint, our analysis covers a broad array of six datasets derived from macroeconomics, microeconomics, and finance. Five of these datasets are in line with those used by Giannone et al. (2022), and one is sourced from Gu et al. (2020). Our finance examples delve into predicting market returns using financial and economic indicators, as well as firm-level return prediction based on their specific characteristics. In the macroeconomic context, we examine time-series predictions of industrial production using macroeconomic indicators, and a cross-country GDP growth prediction, utilizing socio-economic, institutional, and geographical factors. Our microeconomic case studies focus on crime rate predictions and pro-plaintiff appellate decisions in takings law rulings.

The relevance of weak signals in datasets is contingent on the choice of benchmark models. For instance, when compared to a constant benchmark model, weak signals are revealed in four out of six datasets. Further benchmarking against covariates informed by economic theory reveals weak signals across all datasets, making them particularly well-suited for the application of our asymptotic theory. Drawing from their empirical analysis of these datasets, Giannone et al. (2022) argue that sparsity may be an illusion, as optimal predictive models often rely on a large number of covariates. Our collective theoretical and empirical evidence points to signal weakness as a key factor in the underperformance of Lasso. As our simulations suggest, even in cases where the majority of signals are spurious — that is, their coefficients are zero in the true DGP — Ridge may still outperform Lasso if the non-spurious signals are weak. This comparative analysis of their performances does not necessarily offer insights into the nature of the DGP itself. In light of these findings, we recommend a cautious
approach to employing Lasso in economic and financial settings. Despite its popularity as a modern counterpart to OLS, Lasso’s effectiveness may be compromised in scenarios characterized by weak signals.

Our paper is closely related to the literature examining the theoretical performances of Ridge and Lasso, with two main threads of results being primarily relevant. The first strand is centered on the error-bound analysis. For Ridge, Hoerl and Kennard (1970) provided an explicit formula for the mean squared prediction error. Based on this, it becomes evident that given a design matrix with bounded singular values, the prediction error vanishes at a rate of $p/n$, with $p$ denoting the number of covariates and $n$ indicating the sample size. More generally, the magnitude of the prediction error is intrinsically tied to the eigen-structure of the design matrix. On the Lasso front, the prediction error vanishes if $s \log p/n \to 0$, where $s$ represents the number of non-zero parameters. This result has been well-documented in a series of studies, including Zou (2006), Zhao and Yu (2006), Zhang and Huang (2008), Bickel et al. (2009), among others. Nevertheless, the unique asymptotic setting we delve into situates itself in a regime where the aforementioned error bounds are insufficient in differentiating the performance of these estimators with the zero estimator, given that the leading order of their prediction errors is identical. This circumstance necessitates a more granular higher-order analysis of prediction errors.

The second, more recent strand of research determines the precise probability limit of the prediction error for Ridge and Lasso. Bayati and Montanari (2012) was at the forefront, employing approximate message passing algorithm and establishing its link with the Lasso estimator. By examining the limiting evolutionary path of this algorithm, they pinpointed the limit of the error. In a different approach, Thrampoulidis et al. (2015) harnessed the Convex Gaussian Minimax Theory (CGMT) to streamline the original optimization problem of Lasso, facilitating the derivation of the precise error. Concurrently, Ridge regression underwent analogous progress, with Dicker (2016) providing insights into the precise errors of Ridge estimators. However, the pursuit of precise error analysis often demands more stringent parametric assumptions. Notably, these studies impose the assumption that the design matrix’s elements are independently Gaussian-distributed. Dobriban and Wager (2018) expanded Dicker (2016)’s analysis of Ridge estimator to accommodate dependent covariates and non-Gaussian predictors, leveraging the universality results in random matrix theory.

The technique of precise error analysis has proven invaluable for gaining deeper insights into the properties of various machine learning techniques. For instance, Liang and Sur (2022) employed CGMT to probe the properties of both minimum $\ell_1$-norm interpolation

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and boosting in linear models. Miolane and Montanari (2021) illuminated the properties of cross-validation for Lasso. Further, Hastie et al. (2022) delved into minimum $\ell_2$-norm interpolation, elucidating the double-descent pattern observed in neural networks and highlighting the potential advantages of overparametrization. On variable selection, Su et al. (2017) investigated the false discovery rate of the Lasso path, while Wang et al. (2020) presented a comparative analysis of the variable selection properties of bridge estimators.

Several studies, including those by Donoho and Jin (2004), Hall and Jin (2010), and Jin and Ke (2016), have explored variable selection in the context of rare and weak signals. Our primary focus, however, is on prediction. Merely identifying variables with non-zero coefficients does not necessarily translate to improved out-of-sample prediction, as their inclusion is also contingent on the bias-variance tradeoff. Additionally, predictors without variable selection properties, like Ridge, might not underperform compared to those that possess such properties, such as Lasso. Moreover, in many asymptotic settings, the task of identifying variables with non-zero coefficients often proves infeasible. Our paper addresses precisely this scenario, albeit with a focus on the aspects of prediction.

In econometrics, the challenges posed by “weakness” are a pivotal area of research. Within the literature on causal inference, the issue of weak instrumental variables has garnered significant attention. A series of robust estimation and inference methodologies have been proposed by Staiger and Stock (1997), Stock and Wright (2000), Andrews and Cheng (2012), Andrews and Mikusheva (2016), to name a few. In parallel, the literature on factor models has evolved with diverse strategies to tackle the challenge of weak factors. Notably, Onatski (2010) introduced tests for determining the number of factors. The task of estimating factor space, in different contexts, has been handled by Freyaldenhoven (2022), Uematsu and Yamagata (2022), Bai and Ng (2023), and Giglio et al. (2023).

This paper is organized as follows. Section 2 presents the main theoretical results regarding Ridge and Lasso regressions. Section 3 conducts simulations to illuminate our theoretical predictions while also expanding the analysis to assess the performance of advanced machine learning methods under weak signals. Lastly, Section 4 provides empirical results supporting the practical relevance of our theoretical results.

**Notation:** For any $x \in \mathbb{R}$, we refer to $\max(x, 0)$ as $x_+$. For any vector $x$, $\|x\|_1$, $\|x\|$ and $\|x\|_\infty$ represent its $\ell_1$, $\ell_2$ and $\ell_\infty$ norms, respectively. For a matrix $A$, we use $\|A\|$ and $\|A\|_F$ to denote the spectral norm (or $\ell_2$ norm), and the Frobenius norm of a matrix $A$, that is, $\sqrt{\lambda_{\max}(A^\top A)}$, and $\sqrt{\text{Tr}(A^\top A)}$, respectively. In the case where $A$ is a $p \times p$ matrix, $\lambda_i(A)$ denotes its $i$-th largest eigenvalue, for $1 \leq i \leq p$. Furthermore, we define its spectral...
distribution as

\[ F^A(x) = \frac{1}{p} \sum_{k=1}^{p} 1_{\lambda_k(A) \leq x}. \]

We use the notation \( x_n \lesssim y_n \) when there exists a constant \( C \) such that \( x_n \leq C y_n \) holds for sufficiently large \( n \). Similarly, we use \( x_n \lesssim_P y_n \) to denote \( x_n = O_P(y_n) \). If \( x_n \lesssim y_n \) and \( y_n \lesssim x_n \), we write \( x_n \asymp y_n \) for short. Similarly, we use \( x_n \asymp_P y_n \) if \( x_n \lesssim_P y_n \) and \( y_n \lesssim_P x_n \).

2 Theoretical Results

2.1 Model Setup

We start with the following linear regression model:

\[ y = X\beta_0 + \varepsilon, \tag{1} \]

where \( y \in \mathbb{R}^n \), \( X \in \mathbb{R}^{n \times p} \), \( \beta_0 \in \mathbb{R}^p \) and \( \varepsilon \in \mathbb{R}^n \).

Our objective is to investigate the performance of machine learning techniques when dealing with weak signals. To accomplish this, we focus on a high-dimensional regression setting characterized by an increasing number of predictors, that is, \( p \rightarrow \infty \). In such a context, regularization techniques become not just relevant but often necessary due to the challenges posed by the curse of dimensionality.

Moreover, our specific focus is on situations where the signals are weak, characterized by the condition: \( \|\beta_0\| \asymp \tau \rightarrow 0 \). The choice to use \( \|\beta_0\| \) as the metric for characterizing weak signals is due to its close relationship with the widely-adopted \( R^2 \) metric in regression analysis, which provides a familiar and intuitive understanding of signal strength.

Our investigation then delves into an asymptotic analysis within a regime broadly characterized by these two conditions. The exact conditions \( p \), \( \tau \), and the sample size \( n \) satisfy will be provided in detail once we introduce the baseline estimator.

Now, we proceed to present the assumptions governing the DGP of \( X \):

**Assumption 1.** The covariates \( X \in \mathbb{R}^{n \times p} \) are generated as \( X = \Sigma_1^{1/2} Z \Sigma_2^{1/2} \) for an \( n \times p \) matrix \( Z \) with i.i.d. standard Gaussian entries, deterministic \( n \times n \) and \( p \times p \) positive definite matrices \( \Sigma_1 \) and \( \Sigma_2 \).\(^2\) In addition, there exist positive constants \( c_1, C_1, c_2, C_2 \) such that \( c_1 \leq \lambda_i(\Sigma_1) \leq C_1, i = 1, 2, \ldots, n \) and \( c_2 \leq \lambda_i(\Sigma_2) \leq C_2, i = 1, 2, \ldots, p \).

\(^2\)We are also able to accommodate random \( \Sigma_1 \) and \( \Sigma_2 \), with an additional assumption that their entries are mutually independent and also independent of \( Z \).
This assumption effectively accommodates both time series and cross-sectional dependence among the covariates in $X$, with $\Sigma_1$ capturing both heteroskedasticity and autocorrelations, while $\Sigma_2$ characterizes cross-sectional correlations. The constraints placed on the eigenvalues of $\Sigma_1$ and $\Sigma_2$ serve a dual purpose. First, the upper bounds on these eigenvalues eliminate strong time series and cross-sectional dependencies within $X$. Second, the lower bounds prevent multicollinearity and the scenario where an observation at a particular time point is linearly dependent on observations from other times.

While the Gaussian assumption for $X$ is integral to our use of Gordon’s inequality (Gordon (1988)) for Gaussian processes in the proof, it does raise concerns regarding the robustness of our findings when this assumption is not met in real-world scenarios. Our simulation results indicate that the Gaussian assumption appears non-essential and our asymptotic theory approximates finite sample behavior even with relatively small sizes, typically a few hundred observations. This observation aligns with similar findings in random matrix theory, where asymptotic properties initially derived for Gaussian ensembles were subsequently shown to extend to a wider spectrum of random matrices — a phenomenon referred to as the universality, as also noted by Bayati and Montanari (2012). In fact, in cases where $\Sigma_1$ is the identity matrix, Dobriban and Wager (2018) demonstrate the feasibility of conducting precise error analysis using random matrix theory, thus bypassing the Gaussian assumption. However, their technique appears to be only applicable to the Ridge estimator. Our aim is to conduct a comparative analysis of both Ridge and Lasso under a unified framework.

Next, we specify the assumptions regarding $\varepsilon$. Similar to the case of $X$, we introduce $\Sigma_\varepsilon$ to account for heteroskedasticity and autocorrelations in the noise.

**Assumption 2.** Let $\varepsilon = \Sigma_\varepsilon^{1/2}z$, where $z$ comprises i.i.d. variables with mean zero, variance one and finite fourth moment and $\Sigma_\varepsilon$ is a positive semidefinite matrix satisfying $c_\varepsilon \leq \lambda_i(\Sigma_\varepsilon) \leq C_\varepsilon$, $i = 1, 2, \ldots, n$, for some fixed positive constants $c_\varepsilon$ and $C_\varepsilon$.

If $\Sigma_\varepsilon$ is a diagonal matrix, its spectral norm is evidently bounded under the condition that every element along the diagonal is bounded, i.e., that $\varepsilon$ has finite variances. In the appendix, we further establish that even if the noise follows a stationary process characterized by exponentially decaying autocorrelations, the spectral norm of $\Sigma_\varepsilon$ remains bounded.

Under Assumptions 1 and 2, it follows that $\|X\beta_0\| \approx_p \sqrt{n} \|eta_0\|$ and $\|\varepsilon\| \approx_p \sqrt{n}$. This indicates that the magnitude of each entry in matrix $X$ and the error term $\varepsilon$ neither explode nor vanish asymptotically. Consequently, the magnitude of the signal-to-noise ratio (or prediction $R^2$) is entirely dictated by $\|\beta_0\|$. Next, we impose an assumption that governs the properties of a large number of parameters collected in $\beta_0$: 
Assumption 3. The vector $b_0 = \sqrt{p\tau^{-1}}\beta_0$ comprises i.i.d. random variables, each following a prior probability distribution $F$ belonging to the class $\mathcal{F}$. The class $\mathcal{F}$ is defined by sub-exponential distributions with a mean of zero and a variance denoted as $\sigma^2_\beta$.

Without loss of generality, we use the term $\sqrt{p\tau^{-1}}$ as the normalization factor, ensuring that $\|\beta_0\|^2 \asymp_{\mathcal{P}} \tau$. This choice of normalization facilitates a clearer interpretation of our results. While the i.i.d. assumption may seem strong, it offers greater transparency by simplifying more complex technical assumptions necessary to derive essential probability bounds. In particular, this assumption allows for important classes of models, such as a spike-and-slab prior for $b_0$, extensively studied by Giannone et al. (2022) to examine the empirical relevance of sparsity in economic datasets. Specifically, each element of $b_0$ follows a mixed distribution: $q\psi_0 + (1-q)\psi_1$, where $q \in [0, 1]$, $\psi_0$ and $\psi_1$ are the spike and slab components of the prior, respectively. These components may assume the form of Gaussian distributions, as suggested by George and McCulloch (1993), or Laplace distributions, as explored in Rovcková and George (2018). This spike-and-slab model will also be utilized in our simulations to test our theoretical predictions.

The underlying assumptions that justify Ridge and Lasso are notably distinct, particularly in the context of error-bound analysis. For instance, the analysis of Lasso often requires the approximate sparsity condition and the restricted eigenvalue condition (see, e.g., Belloni et al. (2013b) and Bickel et al. (2009)). On the other hand, the convergence rate of Ridge’s prediction error requires intricate conditions on the eigenvalue structure of the design matrix, as discussed in Tsigler and Bartlett (2023). In contrast, our analysis here compares the asymptotic properties of different estimators within a common framework. Central to our analysis is the consideration of the probability limit of the prediction errors, which necessitates assuming a (prior) probability distribution on the coefficients. This choice is in line with common practices in this literature on precise errors, which also connects our analysis of prediction error with Bayes risk.

2.2 Estimators

We now turn our attention to the discussion of the estimators. In scenarios involving weak signals, characterized by $\|\beta_0\| \to 0$, a straightforward and natural baseline estimator emerges, that is, the naive zero estimator. This estimator is clearly consistent in terms of the $\ell_2$-loss of the estimation error, because the error reduces to $\|\beta_0\|$, which vanishes in this context.

The zero estimator functions as a passive baseline, serving as a benchmark for a scenario where no learning occurs. To surpass the performance of the zero estimator, any alternative
estimator must harness some of the available weak signals. This indicates the alternative estimator’s capability to successfully identify and leverage predictive signals, even when they are weak. Therefore, to address the earlier question of whether machines can learn weak signals, we only need to compare the machine learning method’s performance with that of the zero estimator. Only if they can do so can they outperform the naive zero estimator.

In our study, we consider Ridge and Lasso as contenders that leverage machine learning techniques. These methods are widely used benchmarks in practice, owing to their simplicity and universality. An in-depth analysis of these estimators provides valuable insights into their specific regularization techniques, which can be extended to more advanced models.

The Ridge estimator, denoted as $\hat{\beta}_r$, is the solution to the following optimization problem:

$$
\hat{\beta}_r(\lambda_n) := \arg\min_{\beta} \frac{1}{n} \| y - X\beta \|^2 + \frac{p\lambda_n}{n} \| \beta \|^2,
$$

where $\lambda_n$ is its tuning parameter governing the strength of the regularization. In contrast, the Lasso estimator, denoted as $\hat{\beta}_l$, is defined as:

$$
\hat{\beta}_l(\lambda_n) := \arg\min_{\beta} \frac{1}{n} \| y - X\beta \|^2 + \frac{\lambda_n}{\sqrt{n}} \| \beta \|_1.
$$

By convention, and without loss of generality, the terms involving penalties are typically scaled by $p/n$ in the case of the Ridge estimator and by $1/\sqrt{n}$ in the case of Lasso.

In addition to Ridge and Lasso, our theoretical results also encompass the ordinary least squares (OLS) estimator and the Ridgeless estimator, both of which correspond to special cases of Ridge when the tuning parameter $\lambda_n$ is set to zero. When $p \leq n$, the least squares problem yields a unique solution, which is the OLS estimator. However, when $p > n$, the least squares problem has an infinite number of solutions. Among these solutions, the Ridgeless estimator can be regarded as a minimum-norm interpolating linear predictor, aiming to minimize the $\ell_2$-norm of $\beta$:

$$
\hat{\beta}_r(0) = \arg\min_{\beta} \| \beta \|, \quad \text{s.t.} \quad X\beta = y,
$$

as noted by Bartlett et al. (2020). It is also possible to explore other interpolators, such as the minimum $\ell_1$-norm interpolator studied by Liang and Sur (2022). Future research might extend our analysis to other penalized linear estimators, such as Elastic Net, as introduced by Zou and Hastie (2005), or SCAD by Fan and Li (2001).
2.3 Bayes Risk

With $\beta$ estimated, it is straightforward to construct corresponding linear predictors. Now, we proceed to define the metric by which we assess various predictors. For any predictor, our primary interest is its Bayes prediction risk. This risk is related to the expected squared prediction error evaluated at a new, independent data point $(x_{\text{new}}, y_{\text{new}})$. In the case of a linear model, we can write the prediction error explicitly as:

$$
\mathbb{E}_F(y_{\text{new}} - \hat{y}_{\text{new}})^2 = \sigma^2 + \mathbb{E}_F\left\{\mathbb{E}\left[(x_{\text{new}})^\top(\hat{\beta} - \beta_0)^2|X, y, \beta_0]\right]\right\} = \sigma^2 + \mathbb{E}_F\|\Sigma^{1/2}_{\text{new}}(\hat{\beta} - \beta_0)\|^2,
$$

(5)

where the subscript in the expectation operator $\mathbb{E}_F(\cdot)$ emphasizes the fact that the expectation is taken with respect to the prior distribution of $b_0 = \sqrt{p\tau^{-1}}\beta_0$. Given that $\sigma^2$ does not depend on the estimator, it is the second term that governs the relative predictive performance of different estimators. Barring $\|\Sigma_2\|$, the prediction risk is closely tied to the estimation error of $\hat{\beta}$, i.e., $\|\hat{\beta} - \beta_0\|$.

Formally, we define the Bayes prediction risk associated with an estimator $\hat{\beta}$ as

$$
\mathcal{R}(\hat{\beta}, F) := \mathbb{E}_F\|\Sigma^{1/2}_{\text{new}}(\hat{\beta} - \beta_0)\|^2.
$$

The predictor that minimizes this Bayes risk is termed the Bayes predictor. In our framework, it is straightforward to show (see Chapter 4 of Berger (1985)) that the Bayes predictor corresponds with the predictor that is derived from the posterior mean of $\beta_0$, which is represented as $\mathbb{E}_F(\beta_0|X, y)$. We denote its Bayes risk as $\mathcal{R}(F)$.

Under strong signals, i.e., $\tau = 1$, and suppose that $\Sigma_1, \Sigma_2$ and $\Sigma_\varepsilon$ are identity matrices, $p/n \to c_0 \in \mathbb{R}^+$, significant progress has been made in understanding the asymptotic behavior of Bayes risk. For Ridge regression, notable studies by Dicker (2016) and Dobriban and Wager (2018) have derived the asymptotic limit of the Bayes risk.

3There exists an extensive body of literature focused on empirical Bayes methods, which explores feasible approaches for implementing $\mathbb{E}_F(\beta_0|X, y)$, in cases where $F$ is unknown, see, e.g., Robbins (1964), Efron (2012), Brown and Greenshtein (2009), and Jiang and Zhang (2009).

4The exact form of the limit is given by

$$
\lim_{n \to \infty} \mathcal{R}(\hat{\beta}_r(\lambda_n), F) = c_0m(-\lambda, c_0) + \lambda(\lambda c_0^2 - c_0)m'(-\lambda, c_0),
$$

(6)

where $\lambda = \lim_{n \to \infty} c_0\lambda_n$ and

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of Lasso, several studies, such as those conducted by Bayati and Montanari (2012) and Thrampoulidis et al. (2018), have established its asymptotic Bayes risk limit.\footnote{The limit in the case of Lasso can be explicitly written as follows: $\lim_{n \to \infty} R(\hat{\beta}_l(\lambda_n), F) = (\alpha^*)^2$, where}

Figure 2: Comparison of Prediction Errors: Optimal Ridge and Lasso vs. the Zero Estimator

Note: The left panel illustrates the ratio of prediction error between the optimal Ridge and the baseline zero estimator. Conversely, the right panel presents a similar comparison for the optimal Lasso estimator against the zero estimator. Both axes, $y$ and $x$, depict a range of $p/n$ ratios and $\|\beta_0\|^2$ values, corresponding to data generated in accordance with the model described in (1), with $\Sigma_1$, $\Sigma_2$, and $\Sigma_\varepsilon$ in Assumptions 1 and 2 set as identity matrices. In this context of strong signals, the prediction errors for both optimal Ridge and Lasso are calculated using tuning parameters that are optimally selected to minimize the expected prediction errors’ probability limits, as given by (6) and (7).

In Figure 2, we present two heatmaps, with the $y$ and $x$ axes representing various values of $p/n$ and $\|\beta_0\|^2 = \tau \sigma^2_\beta$. The left heatmap illustrates the ratio of Bayes risk between optimal Ridge and the zero estimator, while the right heatmap represents the ratio of optimal Lasso against the zero estimator. For both Ridge and Lasso, their optimal tuning parameters are selected by minimizing the probability limits of their Bayes risk given by (6) and (7),

$$m(-\lambda, c_0) = \frac{-(1 - c_0 + \lambda) + \sqrt{(1 - c_0 + \lambda)^2 + 4c_0\lambda}}{2c_0\lambda}.$$
respectively. A prediction error ratio below 1 within these visualizations suggests that the zero estimator is outperformed.

The heatmaps, as anticipated, clearly demonstrate that both Ridge and Lasso estimators surpass the performance of the zero estimator. This superiority is particularly pronounced in scenarios involving strong signals and relatively lower dimensions. Notably, the disparity between these estimators becomes less pronounced as the norm of $\|\beta_0\|$ approaches zero and the ratio $p/n$ increases, indicating a shift towards scenarios characterized by weaker signals. The existing result on precise error analysis is primarily built upon the assumption of strong signals, where $\tau = 1$. To discern the performance of various estimators under weak signal conditions, a more intricate analysis in the limiting case ($\tau \to 0$ and $p/n \to \infty$) is necessary.

### 2.4 Zero’s Optimality and Relative Prediction Error

Figure 2 also indicates that our attention should be directed towards a regime where the zero estimator exhibits meaningful competitiveness. Otherwise, we may question the appropriateness of our definition of “weak” signals if some machine learning approaches can obviously outperform it.

One might be tempted to define “weak signals” as instances where the signal strength falls below a certain “detection boundary,” thereby becoming indiscernible through hypothesis testing. The “detection boundary” in this scenario represents the threshold level of signal strength at which statistical tests can reliably discern the presence of a signal amidst noise or other influencing factors. Relevant tests include Ingster et al. (2010), Cui et al. (2018), and Li et al. (2020). However, this boundary generally hinges on the chosen alternative and the maintained hypotheses, presenting a challenge in establishing a unified benchmark. More importantly, our primary focus is on prediction, rather than signal detection. This distinction is key because, even when signals are undetectable by hypothesis testing, their collective contribution to prediction can still outperform the zero predictor. The zero predictor also serves as a natural benchmark for demonstrating the capacity of machine learning to identify and utilize weak signals.

To motivate our concept of weak signals, we analyze a regime where the zero estimator achieves certain notion of optimality, indicated by its Bayes risk being identical to that of the Bayes predictor. This scenario is delineated more precisely by the assumption below:

**Assumption 4.** $\tau \to 0$, $p/n \to c_0 \in (0, \infty]$, $\tau p(\log p)^4/n \to 0$, and $n\tau p^{-2/3}(\log p)^{-4} \to \infty$.

Assumption 4 covers a wide spectrum of signal strengths and counts, while simultaneously imposing constraints to prevent an excessively large $p/n$ ratio and too rapid vanishing of $\tau$. 

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These constraints can be visually represented by a specific area around the positive section of the y-axis in Figure 2. The latter two constraints effectively exclude areas near the origin and positive infinity from our analysis, because in these regions, the relative performance of different estimators cannot be deterministically ascertained.

To facilitate our discussion of the optimal estimator, we refer to the definition provided by Robbins (1964).

**Definition 1.** We say $\hat{\beta}$ is asymptotically optimal relative to $F$, if it satisfies

$$\lim_{n \to \infty} \frac{R(\hat{\beta}, F)}{R(F)} = 1.\footnote{The definition of asymptotic optimality is provided in terms of a ratio to accommodate more general scenarios where $R(F)$ varies with the sample size $n$.}

**Theorem 1.** Assume that Assumptions 1–4 hold. Furthermore, assume that the error term $\varepsilon$ in Assumption 2 follows a Gaussian distribution.\footnote{The Gaussian assumption on $\varepsilon$ is only used to facilitate considerations of optimality, which is a standard assumption in the empirical Bayes literature, e.g., Jiang and Zhang (2009). While this assumption motivates our characterization of weak signal regimes, it is not utilized in follow-up analysis of the estimators.} Under these conditions, the zero estimator is asymptotically optimal relative to any distribution $F \in F$.

This theorem suggests that regardless of the unknown prior distribution, the prediction risk associated with the zero estimator is asymptotically identical to that of the Bayes predictor. Notably, while the Bayes predictor requires knowledge of the prior distribution and is thus infeasible in many practical scenarios, the zero estimator achieves the same level of prediction risk without requiring such information. From this perspective, the zero estimator is both feasible and optimal.

It is noteworthy that the zero estimator can be considered as a particular case of both Ridge and Lasso estimators when a sufficiently large tuning parameter is chosen. Given this perspective, and in accordance with the insights of Theorem 1, the relative Bayes risk of the optimal Ridge and Lasso estimators, in comparison to the zero estimator, is expected to asymptotically approach one. This result suggests that under conditions of weak signals, merely comparing their Bayes risk ratios may not be an effective approach to tell any differences among these estimators.

As such, we shift our attention to the relative prediction error between any estimator $\hat{\beta}$ and the zero estimator, defined as follows, in absolute difference rather than their ratio. To ensure a meaningful scale in the limit, we multiply the relative error by $pn^{-1}\tau^{-2}$, and adopt
the following metric for comparison:

\[ \Delta(\hat{\beta}) = pn^{-1}\tau^{-2}(\|\Sigma_2^{1/2}(\hat{\beta} - \beta_0)\|^2 - \|\Sigma_2^{1/2}\beta_0\|^2). \]  

(8)

\(\Delta(\hat{\beta})\) magnifies the relative prediction error, a measure that predominantly depends on higher order differences for estimators we consider. Based on this definition, if \(\Delta(\hat{\beta}) > 0\) holds with probability approaching one, it indicates that the estimator \(\hat{\beta}\) exhibits inferior prediction performance compared to zero. Conversely, if \(\Delta(\hat{\beta}) < 0\) holds with probability approaching one, it implies that the estimator \(\hat{\beta}\) outperforms zero.

Before we proceed to present our main results in the following section, we need to provide technical conditions governing the limiting behavior of \(\Sigma_1, \Sigma_2,\) and \(\Sigma_\varepsilon\):

**Assumption 5.** For matrices \(\Sigma_1, \Sigma_2\) and \(\Sigma_\varepsilon\):

\[
\frac{1}{n} \text{Tr}(\Sigma_1) = 1 + O(n^{-1/2}), \quad \frac{1}{p} \text{Tr}(\Sigma_2) = \sigma_x^2 + O(p^{-1/2}), \quad \frac{1}{n} \text{Tr}(\Sigma_\varepsilon) = \sigma_\varepsilon^2 + O(n^{-1/2}).
\]

Additionally, there exist constants \(\theta_1\) to \(\theta_4\) such that

\[
\frac{1}{n} \text{Tr}(\Sigma_\varepsilon \Sigma_1) = \sigma_\varepsilon^2 \theta_1 + o(n\tau/p),
\]

\[
\frac{1}{p} \text{Tr}(\Sigma_2^2) = \sigma_x^4 \theta_2 + o(1), \quad \frac{1}{n} \text{Tr}(\Sigma_\varepsilon \Sigma_1^2) = \sigma_\varepsilon^2 \theta_3 + o(n/p), \quad \frac{1}{n} \text{Tr}(\Sigma_1^2) = \theta_4 + o(n/p).
\]

As \(\Sigma_1, \Sigma_2,\) and \(\Sigma_\varepsilon\) are positive definite, all of these constants \(\theta_i\), where \(i = 1, 2, 3,\) and 4, are positive. The condition concerning \(\Sigma_2\) can be verified through a more primitive condition often found in the literature—namely, the existence of the limit of \(\Sigma_2\)’s empirical spectral distribution \(F_{\Sigma_2}(x)\), as assumed by Dobriban and Wager (2018). Regarding the conditions concerning \(\Sigma_1\) and \(\Sigma_\varepsilon\), we establish in the appendix (Lemma ??) that when the time series of covariates and noise are stationary with exponentially decaying correlations, these conditions hold. In situations where all three matrices reduce to identity matrices, which is a common scenario in the literature on precise error analysis, Assumption 5 holds trivially.

### 2.5 Analysis of the Ridge Estimator

In this section, we present the results of the Ridge estimator in the context of weak signals. We begin by presenting the relative error of Ridge for any tuning parameter value:

**Theorem 2.** Assuming that Assumptions 1–5 hold, and setting \(\lambda_n = \tau^{-1}\lambda\), we establish the following convergence result:
\[
\Delta(\hat{\beta}_r(\lambda_n)) \xrightarrow{p} \alpha^* := 2\theta_x\sigma_x^4 \left( \frac{\sigma_x^2\theta_1}{2\lambda^2} - \frac{\sigma_\theta^2}{\lambda} \right).
\]

This theorem yields several important findings. First, by minimizing \(\alpha^*\) with respect to \(\lambda\), we can determine the optimal tuning parameter value: \(\lambda_n^{\text{opt}} = \tau^{-1}\sigma_x^2\theta_1/\sigma_\theta^2\). Furthermore, with the optimal tuning parameter in place, \(\alpha^*\) is negative, indicating that Ridge can effectively learn weak signals when the tuning parameter is chosen appropriately.

Second, when we set \(\lambda \to \infty\) (equivalently, \(\tau\lambda_n \to \infty\)), the value of \(\alpha^*\) converges to zero. This outcome is expected, as the use of a large tuning parameter makes the estimator’s performance increasingly resemble that of the zero estimator. Nevertheless, it is noteworthy that \(\alpha^* \to 0^-\); in other words, as \(\lambda\) increases, the Ridge estimator consistently outperforms the zero estimator until it gradually becomes indistinguishable from it in the limit.

Third, as \(\lambda \to 0\), in which case \(\lambda_n = o(\tau^{-1})\), the corresponding value of \(\alpha^*\) tends to positive infinity. This indicates that Ridge’s performance deteriorates to the point where the Ridgeless estimator (corresponding to \(\lambda = 0\)) is surpassed by the zero estimator. This is a significant departure from the strong signal setup in which Ridgeless can still outperform the zero estimator, as demonstrated by Hastie et al. (2022). It is important to note that the Ridgeless estimator, defined in the form of no regularization (\(\lambda = 0\)), is not completely devoid of regularization. It incorporates implicit regularization by yielding the interpolator that achieves the minimum \(\ell_2\) norm. This inherent form of regularization enables the Ridgeless estimator to effectively control variance, particularly in situations where the number of predictors \(p\) exceeds the sample size \(n\), thus ensuring desirable performance in strong signal scenarios. In contrast, under conditions of weak signals, this implicit regularization is insufficient for effective variance control. This inadequacy results in the estimated \(\|\hat{\beta}\|\) being substantially larger than \(\|\beta_0\|\), leading to the poor performance of \(\hat{\beta}\).

Furthermore, given that the Ridgeless estimator is defined as the interpolator that minimizes \(\|\hat{\beta}\|\), it follows that all linear interpolators, including, for instance, the one that minimizes the \(\ell_1\)-norm, result in even larger values of \(\|\hat{\beta}\|\). Consequently, these interpolators also fail to outperform the zero estimator in contexts with weak signals.

Figure 3 provides an illustrative example of the relationship between the relative error of the Ridge estimator and the tuning parameter \(\lambda\), showcasing the theoretical insights we have discussed. Corollary 1 below summarizes the result on the minimal-norm interpolating estimator:

**Corollary 1.** Under the same assumptions as in Theorem 2, the Ridgeless estimator, defined by (4), satisfies:
Note: In this plot, the black curve represents the probability limit of $\Delta(\hat{\beta}_r(\lambda_n))$, denoted as $\alpha^*$, as a function of the tuning parameter $\lambda$, defined in Theorem 2, in the context of weak signals. To create this plot, we set all parameters ($c_0, \theta_1, \theta_2, \sigma_x, \sigma_\beta, \sigma_\epsilon$) to one for simplicity.

$$\Delta(\hat{\beta}_r(0)) \overset{P}{\rightarrow} \infty.$$ 

Given Ridge regression’s ability to effectively learn weak signals with an appropriately tuned parameter, the data-dependent selection of this parameter becomes crucial. A paradigmatic approach for this purpose is $K$-fold cross-validation (CV).

In cases where the signals are strong, both Liu and Dobriban (2020) and Hastie et al. (2022) have demonstrated the effectiveness of CV for Ridge. Specifically, the cross-validated tuning parameter converges in probability to the optimal value within a pre-specified interval. The fact that this optimal value lies in some known interval simplifies the derivation of the theoretical properties of CV. In scenarios with weak signals, however, the optimal tuning parameter tends to diverge as the sample size increases. The rate of divergence depends on the unknown strength of the weak signal, $\tau$. As we show next, CV remains a valid and useful tool in this case. To narrow our focus to the matter of weak signals without delving into a complicated CV procedure, we consider the case where both $\Sigma_\epsilon$ and $\Sigma_1$ are identity matrices. This assumption of no temporal dependence in the data facilitates a more straightforward CV procedure for i.i.d. data.

To determine the optimal tuning parameter using $K$-fold CV, denoted as $\hat{\lambda}^{K-CV}$, we begin by partitioning the rows of the design matrix $X$ into $K$ distinct subsets, labeled as $X(1), \cdots, X(K)$. For each index $i \in \{1, \cdots, K\}$, we define $X_{(-i)}$ as the submatrix of
$X$ obtained by excluding the rows corresponding to $X(i)$. Similarly, we have associated subvectors $y(i), \varepsilon(i)$, as well as $y(-i), \varepsilon(-i)$. We next define $\hat{\beta}_r^i(\lambda_n)$ for each $\lambda_n$ as the solution to the Ridge optimization problem for each index $i = 1, \cdots, K$:

$$\hat{\beta}_r^i(\lambda_n) = \arg\min_{\beta} \left\{ \frac{1}{n} \|y(-i) - X(-i)\beta\|^2 + \frac{p\lambda_n}{n} \|\beta\|^2 \right\}.$$  

Consequently, the tuning parameter selected by $K$-fold CV is given by

$$\hat{\lambda}_n^{K-CV} = \arg\min_{\lambda_n \in [\epsilon, \infty)} \frac{1}{n} \sum_{i=1}^{K} \|y(i) - X(i)\hat{\beta}_r^i(\lambda_n)\|^2,$$

where $\epsilon > 0$ is an arbitrary small constant. The following theorem provides justification for the validity of this CV procedure in the context of weak signals:

**Theorem 3.** Under the same assumptions as in Theorem 2, if we also assume that $\Sigma_1 = I$, $\Sigma_\varepsilon = \sigma_\varepsilon^2 I$ and that $\varepsilon$ follows a sub-exponential distribution, then we can establish that:

$$\tau \hat{\lambda}_n^{K-CV} \xrightarrow{P} \lambda_{opt} = \sigma_\varepsilon^2 / \sigma_\beta^2.$$

This theorem justifies the use of $\hat{\lambda}_n^{K-CV}$ as an approximation for the optimal tuning parameter $\lambda_{opt} = \lambda_{opt} / \tau$ ($\theta_1 = 1$ in this case) for Ridge. Importantly, this result does not require prior knowledge of $\tau$, making the CV approach directly applicable in practical scenarios. With our analysis of Ridge concluded, we will now turn our attention to Lasso.

### 2.6 Analysis of the Lasso Estimator

Unlike Ridge, the analysis of Lasso is more intricate, primarily because the Lasso estimator lacks a closed-form formula. In the special case where $\Sigma_1, \Sigma_2, \Sigma_\varepsilon$ are identity matrices, several studies, including Bayati and Montanari (2012) and Thrampoulidis et al. (2018), have established Lasso’s precise error given by (7). Additionally, based on (7), Wang et al. (2020) conducted a small-signal Taylor expansion of $\alpha^*$ with respect to $\sigma_\beta^2$, which affects $\alpha^*$ through the prior distribution $F$. They concluded that the optimal Lasso estimator fails to outperform zero.\(^8\) In the general case we consider, pinpointing the exact precise error appears a daunting task. Instead, we seek probability bounds that allow us to characterize the location of the

\(^8\)The Taylor expansion approach may lack rigor in that $\alpha^*$ itself is already the outcome of a probability limit, and expanding it around a certain parameter could raise concerns about the validity of switching the order of limits.
This turns out sufficient for us to conclude that Lasso cannot outperform zero for all values of its tuning parameter in the context of weak signals. The next theorem summarizes our main findings:

**Theorem 4.** Assume that Assumptions 1–5 are satisfied and the tuning parameter $λ_n$ is chosen such that the following equation holds for some $C_λ > 0$:

$$pn^{-2}τ^{-2}E_{U \sim N(0, Σ_2)} \left\{(2σ_ε \sqrt{θ_1} |U| - λ_n)_+ \right\}^2 = C_λ.$$  \(9\)

Then, with probability approaching one, we have $c_α ≤ Δ(\hat{β}_l(λ_n)) ≤ C_α$, where $c_α$ and $C_α$ are the solutions to the following equation in terms of $x$:

$$x - \sqrt{\frac{2C_λ}{c_2}} x = -\frac{C_λ}{100C_2},$$  \(10\)

where $c_2$ and $C_2$ are constants defined in Assumption 1.

Equation (9) implicitly determines the rate at which $λ_n$ diverges to infinity. For any fixed $C_λ > 0$, we can solve for the tuning parameter $λ_n$ from (9), and derive the upper and lower bounds, $C_α$ and $c_α$, from equation (10). Furthermore, equation (10) directly implies that $C_α$ and $c_α$ are non-negative, indicating that Lasso does not outperform the zero estimator for any given tuning parameter value in the context of weak signals.

Moreover, as $λ_n$ approaches zero, $C_λ$ diverges to infinity, leading to a simultaneous divergence of both $c_α$ and $C_α$. This suggests that the Lasso estimator behaves increasingly worse compared to the zero estimator. Conversely, a larger tuning parameter $λ_n$ causes $C_λ$ to converge to zero from the positive side. As a result, both $c_α$ and $C_α$ converge to zero while remaining non-negative. This implies that the performance of the Lasso estimator improves but remains inferior to the zero estimator, until they become equivalent in the limit.

Figure 4 visually represents upper and lower bounds for the relative error of Lasso in comparison to the zero estimator across various tuning parameter values. This figure aids in conveying the theoretical insights we have discussed earlier.

In light of Lasso’s underwhelming performance, it is evident that in situations involving weak signals, the elastic net estimator, which combines $ℓ_1$ and $ℓ_2$ norms of the parameters in its penalty function, is unlikely to outperform Ridge.\(^{10}\)

\(^9\)When applied to a vector, $| \cdot |$ and $(\cdot)_+$ represent element-wise operations.

\(^{10}\)Although a formal justification for this observation can be provided in the setting of $Σ_2 = I$, we omit it here due to space constraints.
Note: In this plot, the black curves represent the lower and upper probability bounds on $\Delta(\hat{\beta}_l(\lambda_n))$, i.e., $c_\alpha$ and $C_\alpha$, as a function of the tuning parameter $\lambda_n$ in the context of weak signals. In this setup, we fix $n = p = 2,000$. We assign the elements of $b_0$ to follow a standard Gaussian distribution and set $\Sigma$ as the identity matrix $I$. As in Figure 3, we set all parameters $(c_0, \theta_1, \theta_2, \sigma_x, \sigma_\beta, \sigma_\varepsilon)$ to one. Finally, we select $\tau = 0.001$, which results in a population $R^2$ around 0.1%.

In contemporary regression analysis, particularly in the context of a large number of covariates, Lasso has gained prominence as a valuable tool, often regarded as the modern counterpart to OLS. Nevertheless, our analysis has uncovered an important caveat. In situations where the signals are weak, our findings indicate that Lasso, regardless of the choice of its tuning parameter, is not a viable option. This revelation holds substantial implications, especially in fields like economics and finance, where large-scale regression analyses are commonplace, and signal-to-noise ratios tend to be low.

Conversely, our results strongly advocate for the use of Ridge regression in scenarios characterized by weak signals. This insight underscores the critical importance of carefully assessing the data’s specific characteristics and the strength of the underlying signals when making decisions regarding the most suitable regularization technique.

2.7 Assessing Signal-to-Noise Ratio

In line with this perspective, we delve into the assessment of the signal-to-noise ratio, a measure that can offer valuable insights into the viability of different machine learning techniques. Our preceding analyses provide initial but indirect guidance in this regard. Specifically, if
Lasso underperforms the zero estimator, it implies a potential issue with the strength of the signal in the data.

A more conventional and direct approach to evaluating the signal-to-noise ratio is through the goodness-of-fit measure known as $R^2$. However, in-sample $R^2$ is prone to overfitting, and as such, out-of-sample $R^2$ is commonly used in machine learning. This metric essentially involves the comparison of mean-squared errors between two predictors. For our specific application, we have chosen to use zero as the reference predictor and define this metric for a given estimator $\hat{\beta}$ as follows:

$$R^2_{oos}(\hat{\beta}) = 1 - \frac{\sum_{i \in OOS}(y_i - X_i\hat{\beta})^2}{\sum_{i \in OOS}y_i^2}, \quad (11)$$

where “OOS” represents the out-of-sample data.

Since a model’s predictive performance hinges on the signal-to-noise ratio, it is reasonable to employ this metric to evaluate the signal-to-noise ratio inherent in the DGP. In situations characterized by strong signals, it is expected that the out-of-sample $R^2$ serves as a consistent estimator for the signal-to-noise ratio as measured by the population $R^2$. This holds true irrespective of the specific estimator $\hat{\beta}$ employed, as long as it is consistent with respect to $\beta_0$, i.e., $\|\hat{\beta} - \beta_0\| = o_P(1)$. However, in situations characterized by weak signals, the outcome depends critically on the choice of estimator beyond the signal-to-noise ratio itself. As an illustration, both Lasso and Ridge are consistent in the sense that their prediction errors asymptotically diminish, yet the out-of-sample $R^2$ for Lasso can turn non-positive, indicating either no improvement or underperformance compared to the zero estimator, as we have shown in Theorem 4.

In the context of weak signals, the next proposition provides a theoretical justification for the relevance of optimal Ridge’s $R^2_{oos}$ in assessing the signal-to-noise ratio in the data.

**Proposition 1.** Under the same assumptions as Theorem 3, and assuming that the out-of-sample data follows the same DGP as the in-sample data, if $n_{oos}p^{-2}n^2\tau^2 \rightarrow \infty$, where $n_{oos}$ is the size of the out-of-sample data, then for the optimal Ridge estimator, it holds that

$$R^2_{oos}(\hat{\beta}_r(\lambda^{opt}_n)) = p^{-1}n\theta_2(R^2)^2 (1 + o_P(1)),$$

where $R^2$ denotes the population $R$-squared, given by $\tau\sigma_x^2\sigma_\beta^2/(\tau\sigma_x^2\sigma_\beta^2 + \sigma_\epsilon^2)$ in this context.

Interestingly, when the size of out-of-sample data is sufficiently large, the out-of-sample $R^2_{oos}$ is approximately proportional to the squared population $R^2$. The specified condition on
\( n_{oos} \) ensures that the estimation error in out-of-sample \( R^2_{oos} \) does not mask the performance differential between the optimal Ridge and the zero estimator.

While \( R^2_{oos} \) does not exactly mirror the population \( R^2 \), this result implies that \( R^2_{oos} \) still serves as an indicator of signal strength in the data. The reason for their discrepancy is that in the weak signal case, the numerator of the \( R^2_{oos} \)—which reflects the relative prediction error between the two estimators—decreases more rapidly than the numerator of \( R^2 \). Therefore, the numerator of \( R^2_{oos} \) only provides a higher-order characterization of signal strength.

### 2.8 Mixed Signal Strengths and Alternative Benchmarks

In the preceding sections, our analysis primarily focused on scenarios where all signals are weak, leading us to consider the zero estimator as our natural benchmark. This section, however, expands our analysis to include models where potentially strong signals serve as benchmarks. Consider another DGP:

\[
y = W\gamma_0 + X\beta_0 + \varepsilon, \tag{12}
\]

where \( W \in \mathbb{R}^{n \times d} \) represents a predefined set of covariates. These covariates include potentially strong signals and form the basis of the benchmark model. We allow the dimension \( d \) to increase to \( \infty \), however, it does so at a slower rate compared to \( n \), ensuring that OLS of \( y \) against \( W \) remains a viable method for estimation.

In many cases, \( W \) could simply be a vector of ones, allowing us to remain agnostic about the magnitude of the regression’s intercept. In our empirical analysis, \( W \) can be motivated from economic theory, whose impact on the response variable is of central interest. Alternatively, \( W \) can encompass lagged values of \( y \), thereby facilitating the inclusion of temporal dependence in the benchmark model. This setup is particular relevant when using an autoregressive model as a benchmark for forecasting economic variables. Exploring the possibility of a data-driven selection of \( W \) is an intriguing direction for future research, and we plan to delve into this in subsequent studies.

Building on these considerations, our focus now shifts to evaluating and comparing the performance against the OLS benchmark with covariates in \( W \). In this context, the OLS benchmark predictor, \( \hat{y}_{\text{new}}^b \), for a new observation \((w^{\text{new}}, x^{\text{new}})\) is defined as follows:

\[
\hat{y}_{\text{new}}^b = \left( w^{\text{new}} \right)^\top \hat{\gamma}, \quad \text{where} \quad \hat{\gamma} = (W^\top W)^{-1}W^\top y. \tag{13}
\]

The inclusion of \( W \) leads us to explore the following Ridge estimator with regularization
only imposed on coefficients of $X$:

$$
\hat{\beta}(\lambda_n) := \arg\min_{\beta} \left\{ \min_{\gamma} \left( \frac{1}{n} \| y - W\gamma - X\beta \|^2 + \frac{p\lambda_n}{n} \| \beta \|^2 \right) \right\}
= \arg\min_{\beta} \left\{ \frac{1}{n} \| M_Wy - M_WX\beta \|^2 + \frac{p\lambda_n}{n} \| \beta \|^2 \right\},
$$

(14)

where $M_W = I - W(W^TW)^{-1}W^T$. Consequently, the estimator for $\gamma$ is thus given by

$$
\hat{\gamma}(\lambda_n) = (W^TW)^{-1}W^T(y - X\hat{\beta}(\lambda_n)).
$$

(15)

The construction for Lasso is similar. Therefore, utilizing the estimated parameters $(\hat{\beta}(\lambda_n), \hat{\gamma}(\lambda_n))$ we are able to formulate a predictor for $y$ as

$$
y_{new}^* = (w_{new}^*)^T \hat{\gamma}(\lambda_n) + (x_{new}^*)^T \hat{\beta}(\lambda_n) = \hat{y}_{new}^b + (\hat{x}_{new}^*)^T \hat{\beta}(\lambda_n),
$$

(16)

where

$$
\hat{x}_{new}^* = x_{new}^* - X^TW(W^TW)^{-1}w_{new}.
$$

Notably, equation (16) illuminates the role of the second term, $(\hat{x}_{new}^*)^T \hat{\beta}(\lambda_n)$, specifically highlighting the contribution of weak signals in contrast to the OLS benchmark, $\hat{y}_{new}^b$. Moreover, a comparison with the zero-benchmark scenario, previously analyzed, reveals a notable distinction in the modification of the regressor and covariates in equation (14). In this instance, our approach involves a regression of $M_Wy$ against $M_WX$, which intuitively means predicting the residuals of the benchmark model using covariates that have been adjusted to remove the dependence on $W$. While our earlier conclusions are likely still valid, the inclusion of generated variables in regressions brings an additional layer of statistical error that warrants careful examination. The forthcoming theorem will elucidate that this extra error does not compromise our prior conclusion.

Given this context and our previous comparative analysis, we focus on the optimal Ridge estimator in this scenario. This is because the performance of Ridgeless, OLS, or Lasso is unlikely to show improvement with the incorporation of additional estimation error.

**Theorem 5.** Let the design matrix $X$ be generated as $X = W\eta_0 + U$. Assume that the triplet $(U, \beta_0, \varepsilon)$ follows the same distribution as $(X, \beta_0, \varepsilon)$ in Theorem 2. Additionally, the matrix $W$ is independent from $U, \beta_0, \varepsilon$. Each covariate within $W$ is assumed to have a finite second moment. Furthermore, we assume that $d = o(n^2p^{-1}\tau)$, and the eigenvalues of $n^{-1}W^TW$ are
lower bounded by some positive constant. Given these assumptions, the predictor, $\hat{y}^{new}$, as defined in (16) and based on the Ridge estimator from (14) with $\lambda_n = \tau^{-1} \lambda$, and the benchmark predictor $\hat{y}_b^{new}$ from (13), satisfy the following:

$$pn^{-1} \tau^2 \left( \mathbb{E}_F \left[ (\hat{y}^{new} - y^{new})^2 \mid I \right] - \mathbb{E}_F \left[ (\hat{y}_b^{new} - y^{new})^2 \mid I \right] \right) \overset{p}{\rightarrow} \alpha^*, \quad (17)$$

where $I$ denotes the information set generated by $(W, X, y, \gamma_0, \beta_0)$, $\alpha^*$ is defined in Theorem 2, and the tuple $(y^{new}, w^{new}, x^{new})$ satisfies (12).

This result indicates that a Ridge-augmented benchmark model demonstrates superior performance compared to the benchmark model alone. In essence, this suggests that Ridge estimator’s predictive performance retains its superiority over the zero estimator. Notably, the error arising from the initial estimation of the benchmark model does not influence the comparative performance between the Ridge and zero estimators.

3 Monte Carlo Simulations

In this section, we conduct simulation experiments to assess the finite sample performance of our asymptotic theory. We begin by establishing a linear model setup and evaluate the performance of Ridge and Lasso estimators.

3.1 Ridge and Lasso for Linear Models

Now, we provide details of the DGP given by (1) for the first simulation exercise. We set $(\Sigma_1)_{ij} = 2^{-|i-j|}$ for $1 \leq i, j \leq n$. We construct $\Sigma_\epsilon$ as a diagonal matrix with i.i.d. entries sampled from the uniform distribution $U(0.5, 1.5)$. The eigenvalues of $\Sigma_2$ are also simulated from $U(0.5, 1.5)$, and we generate an orthogonal matrix randomly as their corresponding eigenvectors. Together, they constitute $\Sigma_2$. These matrices are generated once and then fixed throughout simulations. By direct calculations, we have $\theta_1 = 1$, $\theta_2 = 13/12$, and $\theta_3 = \theta_4 = 5/3$.

We experiment with two sample sizes, $n = 200$ and $n = 2,000$, while maintaining $p/n = 1$. For each simulated sample, we construct $\beta_0$ as $\sqrt{p^{-1} \tau} b_0$, with $b_0$ drawn from a spike-and-slab distribution: $q\delta_0 + (1 - q)\mathcal{N}(0, (1 - q)^{-1} \sigma^2_\beta)$. Here, $\delta_0$ represents the Dirac delta function, and we set $q = 0.5$ and $\sigma^2_\beta = 1$. As a result, roughly 50% of the coefficients in $\beta_0$ are zero. The error term $\epsilon$ follows a $\mathcal{N}(0, 1)$ distribution. The design matrix $X$ is constructed from $\mathcal{N}(0, 1)$, which are then multiplied by $\Sigma_1^{1/2}$ on the left side and $\Sigma_2^{1/2}$ on the right side. Finally,
we set $\tau = n^{-0.3}/4$, so that the resulting population $R^2$ values range from 4.9% for $n = 200$ to 2.5% for $n = 2,000$. In total, we perform 1,000 Monte Carlo repetitions.

For each simulated sample, we compute the relative prediction error, $\Delta(\hat{\beta}(\lambda_n))$, defined in equation (8). In the case of Ridge regression, we consider three different choices for the tuning parameter, $\lambda_n = \tau^{-1} \lambda$, specifically, $\lambda = 0.5, 1, \text{ and } 2$, where $\lambda = 1$ corresponds to the optimal tuning parameter. The histograms of relative prediction error are presented in Figure 5, with red vertical dashed lines indicating $\alpha^*$ as defined in Theorem 2.

Several noteworthy observations can be made from these histograms. First, across all plots, the probability mass is concentrated around the red vertical line. As the sample size increases from 200 to 2,000, the histograms become increasingly concentrated. This aligns with our theory, which predicts that the relative prediction error converges in probability to the limit $\alpha^*$ as the sample size grows. Second, the value of $\alpha^*$ corresponding to the optimal tuning parameter $\lambda = 1$ is the smallest. This is because the optimal Ridge estimator achieves the smallest prediction error. Moreover, almost all the probability mass corresponding to the optimal Ridge estimator is situated on the negative side of the x-axis, indicating that this estimator outperforms the zero estimator with high probability. Third, when $\lambda = 0.5$, it results in the worst performance, with a large portion of the probability mass on the positive side of zero. In contrast, for $\lambda = 2$, $\alpha^*$ gets closer to zero, and the variance of the relative prediction error decreases. This behavior is due to the increasing amount of penalization, which ultimately drives the estimator towards zero, and in turn, $\alpha^*$ towards zero as well.

Continuing our investigation, we conduct an experiment to analyze $R^2_{oos}$ based on the optimal Ridge. Proposition 1 describes the expected asymptotic behavior of $R^2_{oos}$. To empirically test this, we implement the optimal Ridge, setting $\lambda = 1$, on a training dataset comprising $n = 2,000$ observations. We then generate predictions using a separate test dataset of size $n_{oos}$, and calculate $R^2_{oos}$. We fix $n_{oos} = [\tau^{-3}]$ that meets the theoretical requirement specified in Proposition 1. The comparative analysis between the population $R^2$, the empirically estimated $R^2_{oos}$, and the theoretically derived limit of $R^2_{oos}$ is illustrated in Figure 6. For a clearer visual presentation, we apply a logarithmic transformation to the y-axis. We vary $\tau$ to compare against a range of population $R^2$ values from 0.5% to 10% on the x-axis. The red line represents the average $R^2_{oos}$ over 1,000 Monte Carlo simulations. Additionally, we draw boxplots to describe the distributions of $R^2_{oos}$ across these simulations. The theoretical limit, expressed as $p^{-1} n \theta_2(R^2)^2$, is traced by the blue line, and the green line illustrates the population $R^2$, which would align with a 45-degree line on a standard scale. Notably, in this weak signal setting, the population $R^2$ significantly surpasses the em-
Figure 5: Simulation Results for Ridge in Linear DGPs

$\lambda = 0.5, n = 200$

$\lambda = 1, n = 200$

$\lambda = 2, n = 200$

$\lambda = 0.5, n = 2000$

$\lambda = 1, n = 2000$

$\lambda = 2, n = 2000$

Note: The histograms depict the relative prediction error $\Delta(\hat{\beta}_r(\lambda_n))$ following equation (8) across 1,000 Monte Carlo samples. We consider two different sample sizes ($n = 200$ and $2,000$) and examine three different values of $\lambda_n = \tau^{-1}\lambda$, where $\lambda = 0.5, 1,$ and $2$. Notably, $\lambda = 1$ represents the optimal tuning parameter. The red dashed line indicates the values of $\alpha^*$.

In contrast to the results obtained for Ridge regression, our theoretical framework does not provide a precise error limit for Lasso. Instead, Theorem 4 offers high probability bounds on relative prediction errors. Figure 7 displays histograms of these errors for various tuning parameters and sample sizes, accompanied by two red vertical lines in each plot representing the lower and upper bounds, $c_\alpha$ and $C_\alpha$. These plots yield several interesting findings. First, as the sample size increases, we observe that the probability mass becomes more concentrated and largely falls within the intervals defined by the bounds. Second, regardless of the tuning parameter values, Lasso consistently underperforms the zero estimator in almost all samples when the sample size is large. Third, as the tuning parameter increases (indicated by a decrease in $C_\lambda$), both the lower and upper bounds approach zero. This behavior is a consequence of the increased regularization, which, in turn, steers the estimator closer to zero. In the end, Lasso becomes identical to the zero estimator.
Figure 6: Out-of-Sample $R^2$ for Optimal Ridge in Linear DGPs

Note: The figure presents boxplots showing the distributions of $R^2_{oos}$ for optimal Ridge regression ($\lambda = 1$) over 1,000 Monte Carlo repetitions, with $n = p = 2,000$, $q = 0.5$, and $n_{oos} = 60,000$. We explore a range of population $R^2$ values, from 0.5% to 10% in increments of 0.5% by adjusting $\tau$. The plot features red, blue, and green lines to represent the average $R^2_{oos}$ over Monte Carlo samples, the theoretical limit as given by Proposition 1, and the population $R^2$. In this plot, we employ a logarithmic scale for the y-axis. Without the logarithmic transformation, the green line would align with a 45-degree line. Additionally, the lower boundaries of the boxplots surpass the axis limits in instances where the $R^2_{oos}$ values are negative.

### 3.2 Advanced Machine Learning Methods for Nonlinear Models

Appendix A demonstrates the robustness of our theoretical predictions concerning deviations to model assumptions using Monte Carlo simulations. In this section, we expand our inquiry into the relevance of our theory to nonlinear machine learning methodologies such as RF, GBRT, and NNs, through simulation experiments. While a theoretical analysis of precise error in these algorithms remains challenging, and this part thus involves some speculation, we leverage insights obtained from linear models to elucidate our simulation findings.

We simulate the following DGP, expressed explicitly in element-wise form:

$$y_i = \sum_{j=1}^{p} \beta_{0,j} f(Z_{ij}) + \varepsilon_i, \quad i = 1, \ldots, n,$$

where $y_i$ denotes the $i$th observation of the response variable, $\beta_{0,j}$ represents the coefficient associated with a function $f(\cdot)$ of the predictor variable $Z_{ij}$.

In order to exploit insights from the prior simulation exercise within this new context, we adopt the following procedure for simulating this model: We generate $Z_{ij}$ by applying
Figure 7: Simulation Results for Lasso in Linear DGPs

Note: The histograms depict the relative prediction error $\Delta(\hat{\beta}(\lambda_n))$ following equation (8) across 1,000 Monte Carlo samples. We consider two different sample sizes ($n = 200$ and 2,000) and examine three different values of $C_\lambda$. The two vertical red dashed lines in each figure indicate the values of $c_\alpha$ and $C_\alpha$ that are solutions to (10).

an inverse transform to $X_{ij}$, which was previously simulated in Section 3.1. Specifically, $Z_{ij}$ is defined as $f^{-1}(X_{ij})$, where the design matrix $X$ is constructed using the identical DGP as previously outlined. Additionally, both the coefficients $\beta_0$ and the error term $\varepsilon_i$ follow the same baseline DGP as previously described. This methodology guarantees the replication of the exact simulation results observed when regressing $y$ on $X$. Nevertheless, our primary focus now shifts to predicting $y$ based on nonlinear models of $Z$ without prior knowledge of $f(\cdot)$. The effective signal-to-noise ratio diminishes relative to the linear case due to the added complexity of learning an unknown function $f(\cdot)$. Needless to say, the machine learning models we explore in the subsequent experiments are capable of handling more general DGPs than the one given by Eq. (18).

3.2.1 Simulations with Tree Algorithms

Tree algorithms are widely employed in machine learning, particularly for tackling complex DGPs involving discrete variables, nonlinearities, and intricate covariate interactions. When training a single tree model, it often results in suboptimal performance. To address this limitation, the prevailing strategy is to harness ensemble methods that combine predictions
from multiple trees.

Two prominent regularization techniques used in conjunction with tree models are RF and GBRT. Each of these techniques offers distinct advantages in improving predictive performance. RF employs a technique called bootstrap aggregation, commonly referred to as bagging. This approach involves training numerous trees independently on different bootstrap samples of the dataset. By averaging the predictions from these individual trees, the ensemble model significantly enhances its overall performance. On the other hand, GBRT adopts a boosting approach. Unlike bagging, boosting aims to iteratively improve model performance by fitting residuals from previously trained trees. This iterative process consolidates relatively weak learners into a strong and highly competitive ensemble model.

Since trees are invariant to monotonic transformations, it is only necessary to report their prediction results for the linear DGP, as these will be identical to those of the nonlinear DGP we are considering. However, it is important to recognize that the performance of tree methods may not be as strong as that of Ridge and Lasso. This is partly due to tree methods incurring an additional approximation error as they use piece-wise constant functions to approximate the linear DGP. Therefore, our primary focus should not be on comparing trees with linear models, but rather on evaluating the effectiveness of different ensemble techniques in capturing weak signals and comparing them with the zero predictor.

Our implementation of RF involves two tuning parameters: the depth of each individual tree, which ranges from 5 to 50, and a random selection of features used in each split of the tree. The number of random features depends on the signal-to-noise ratio of the DGP. We select potential values for this tuning parameter within a sufficiently wide grid, while taking into consideration computation expense. The total number of trees in the RF ensemble is fixed at 500. In the case of GBRT, we have three tuning parameters in total, including the depth of the trees, the number of trees, and the learning rate. We adjust the learning rate between 0.001 and 0.5 and vary the depth of each tree from 1 to 6, recognizing that GBRT performs better with shallower trees compared to RF. We observe that for cases where the out-of-sample $R^2$ are 2.5% and 10%, a maximum of 100 trees is sufficient. However, when $R^2$ reaches 40%, the optimal number of trees significantly increases to beyond 10,000. This

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11 Our separate simulation experiments, not included here, also confirm this observation.
12 Specifically, when $R^2$ stands at 2.5%, this parameter spans from 5 to 60; for $R^2$ at 10%, the range extends from 50 to 300; and with $R^2$ at 40%, it falls within the range of 500 to 2,000. The chosen tuning parameter values via CV mostly fall within the middle portion of these ranges, eliminating the need for wider grid searches.
13 Our simulation experiments with GBRT reveal that the selected tuning parameters typically align with the median range of our predefined grid across various scenarios. An exception is noted in the case of the
is expected, because when signal-to-noise ratio is high, the algorithm can identify as many as 2,000 × 0.5 = 1,000 useful covariates, which necessitates a large number of shallow trees. We analyze the benchmark scenario where \( n = p = 2,000, q = 0.5, \) and \( n_{\text{oos}} = 60,000. \) We calibrate \( \tau \) to align with different specified values of the population \( R^2. \) When \( R^2 = 2.5\%, \) both RF and GBRT exhibit underwhelming performance, as illustrated in Figure 8. Neither model surpasses the zero predictor, underscoring their limitations in learning from weak signals. To further assess their capabilities, we increase \( R^2 \) to 10\% and 40\%, exploring their performance under stronger signal-to-noise ratios. At an \( R^2 \) of 10\%, RF demonstrates its potential in learning moderately weak signals, and at 40\%, it markedly outperforms the zero predictor. On the other hand, GBRT faces challenges at the 10\% level, only managing to surpass the zero estimator in most of the simulated samples when \( R^2 \) reaches 40\%. Notably, GBRT demonstrates a higher variance in its prediction error compared to RF. Collectively, this evidence points to RF’s superior ability in learning from weak signals over GBRT. 

A possible explanation for GBRT’s performance pattern could be an inherent \( \ell_1 \)-like regularization in its boosting approach. This hypothesis draws support from the work of Efron et al. (2004), which demonstrates a parallel between boosting and the Lasso path in linear regressions. To substantiate this conjecture, we examine the number of active variables (those with a non-zero importance score) from both tree methods. For RF, the average count of active variables presents a revealing pattern: it is 1999.5 at an \( R^2 \) of 2.5\%, slightly increases to 1999.7 at \( R^2 \) of 10\%, and reaches 2000 at \( R^2 \) of 40\%. This pattern mirrors that of Ridge regression. Conversely, GBRT demonstrates a significantly lower count of active variables, with averages of 28.8, 64.6, and 1736.5 for \( R^2 \) values of 2.5\%, 10\%, and 40\% respectively, aligning more with the variable selection feature of Lasso. Based on our theoretical findings that Ridge outperforms Lasso in weak signal scenarios, we can infer that RF is more adept than GBRT in settings characterized by low signal strengths.

### 3.2.2 Simulations with Neural Networks

Next, we study fully-connected feed-forward NNs, following the stylized architectures as outlined by Gu et al. (2020). For this exploration, we revisit the benchmark case where \( n = p = 2,000 \) and \( q = 0.5. \) Such parameters lead to an input layer in the NN configured with 2,000 neurons. Our specifically chosen architecture features a neural network with a number of trees when \( R^2 = 40\%. \) In this scenario, we observed that reaching the optimal number of trees is hindered by significantly prolonged processing times. Importantly, we also discovered that increasing the number of trees beyond 10,000, up to 20,000, does not yield significant improvements in predictive performance. Consequently, we have decided to limit the total number of trees to 10,000 in this scenario.
Note: The histograms depict the relative prediction error, \( pn^{-1} \tau^{-2} n_{oos}^{-1} \sum_{i \in OOS} (y_i - \hat{y}_i)^2 - y_i^2 \), across 1,000 Monte Carlo samples. We consider RF and GBRT in the case \( n = p = 2,000, q = 0.5, n_{oos} = 60,000 \), and examine three different values of \( R^2 \), which is achieved by adjusting the value of \( \tau \). The red dashed line indicates the y axis.

single hidden layer, which includes 32 neurons.

The training process of these NNs often incorporates a sophisticated mix of optimization and regularization techniques, crucial for enhancing performance. Key methods include stochastic gradient descent (SGD) with Adam (Kingma and Ba (2014)), which expedites the optimization process through an adaptive learning rate. Early stopping, as discussed in Goodfellow et al. (2016), is employed to prevent overfitting by halting training when validation performance starts to decline. Dropout (Srivastava et al. (2014)) is utilized for better generalization, achieved by randomly deactivating neurons. Batch normalization (Ioffe and Szegedy (2015)) aids in stabilizing the training process. Moreover, ensembling over various random seeds is implemented to reduce the variances in model outputs. Furthermore, the integration of \( \ell_1 \) and \( \ell_2 \) penalties with these techniques helps regulate the NN parameters.

To specifically assess the influence of \( \ell_1 \) and \( \ell_2 \) regularization on NN performance, we will minimize potential interference from other factors. Therefore, our implementation will involve using plain SGD as the sole optimization technique, coupled exclusively with either \( \ell_1 \) or \( \ell_2 \) penalties, and deliberately avoiding additional optimization enhancements.\textsuperscript{14} This

\textsuperscript{14}As a result, the training process hinges primarily on two tuning parameters: the learning rate and the regularization parameter. We have set the learning rate to vary between 0.002 and 0.012, a range determined
approach is designed to isolate and clarify the specific contributions of these regularization techniques to NN performance, albeit at the expense of not fully exploiting the NN’s potential. Additionally, given the conceptual similarities between early stopping and shrinkage methods—specifically, early stopping effectively shrinks parameter values towards their initial, smaller magnitudes—we will also conduct a comparative analysis of the effects of early stopping and $\ell_2$-regularization.

To evaluate the performance of NNs, we analyze their behavior with three specific monotonic nonlinear functions: Tangent ($\tan(x)$), Cubic ($x^3$), and Sinh ($(e^x + e^{-x})/2$). The histograms in Figure 9 display the relative prediction errors of NNs when they are applied with various regularization techniques, including early stopping, $\ell_2$ regularization, and $\ell_1$ regularization. It is observed that both $\ell_2$ regularization and early stopping are effective in detecting and leveraging weak signals, evidenced by the majority of the probability mass of their histograms being positioned on the negative side of the y-axis. In contrast, under $\ell_1$ regularization, there seems to be a notable decline in performance.

4 Empirical Analysis of Six Economic Datasets

In this section, we demonstrate the practical relevance of our theoretical insights by applying seven machine learning methods—Ridge, Lasso, OLS/Ridgeless, RF, GBRT, NNs with both $\ell_1$ and $\ell_2$ penalties—across six datasets. These datasets are sourced from three fields: microeconomics, macroeconomics, and finance, with two datasets representing each field. Five of these datasets are similar to those utilized by Giannone et al. (2022), with updates to the latest available data wherever feasible. Additionally, we incorporate an updated dataset from Gu et al. (2020) for our second finance example, which offers a more comprehensive coverage of firm characteristics than the analogous example discussed by Giannone et al. (2022). There is a notable difference between our empirical strategy and that of Giannone et al. (2022), which focuses on estimating a parametric model using a Spike-and-Slab prior within a Bayesian framework. In contrast, our study, aligning more closely with Gu et al. (2020), places a greater emphasis on the comparative analysis of various methods.

At the outset of each empirical exercise, we face a variety of decisions regarding our observations from the validation sample. To efficiently control computational costs, we have adopted a strategy of jointly tuning the learning rate and the number of epochs, while keeping the product of these two factors constant and fixing the batch size in SGD at 200. This method is designed to achieve a balanced compromise, optimizing both the efficiency of the learning process and the stability of the resulting model. On the other hand, the choice of regularization parameter is contingent upon the DGP and the specific regularization technique employed.
Figure 9: Simulation Results for NNs in Nonlinear DGPs

Note: The histograms depict the relative prediction error, $pn^{-1} \tau^{-2} n_{oos}^{-1} \sum_{i \in OOS} (y_i - \hat{y}_i)^2 - y_i^2$, across 1,000 Monte Carlo samples. These histograms pertain to NNs in scenarios where $n = p = 2,000$, $q = 0.5$, $n_{oos} = 60,000$, $\tau = n^{-0.3/4}$, which corresponds to a population $R^2$ of 2.5% with respect to covariates $X_{ij} = f(Z_{ij})$. The focus is on three different regularization techniques: early stopping, $\ell_1$-penalty, and $\ell_2$-penalty, and the experiments encompass three nonlinear models: Tan, Cubic, and Sinh. A red dashed line is used to indicate the y-axis in these histograms.

Implementation strategy. These include defining the in-sample and out-of-sample periods, opting for either a rolling window or an expanding window approach, selecting a cross-validation procedure, and deciding on the normalization of covariates.\textsuperscript{15} We delegate these choices to the frameworks established by Giannone et al. (2022) and Gu et al. (2020), with the intention of minimizing degrees of freedom to enhance the robustness, comparability, and reproducibility of our findings. In the application of each machine learning method, the

\textsuperscript{15}It is crucial to normalize covariates before employing machine learning methods. This step standardizes the scales of covariates, facilitating regularization and enhancing the convergence of optimization algorithms. To avoid forward-looking bias, we ensure that the normalization of covariates is conducted using their respective in-sample mean and standard deviation, thereby maintaining the validity and integrity of our predictive analysis.
selection of an appropriate grid for tuning parameters is essential. This crucial step requires balancing performance optimization with computational efficiency. Finer and wider grids, while potentially enhancing performance, also increase computational demands. Appendix B provides details regarding our model configuration and tuning parameter selection.

Below we present the empirical findings derived from six distinct datasets, each analyzed and reported separately. The primary summary statistics, $R^2_{oos}$, are collected in Table 1. Additionally, we include variable importance plots in Figure 10 as supplementary evidence to decode the performance of different methods. The notion of variable importance is not universally established and varying across different contexts. Our approach diverges from the well-known method associated with RF, originally presented in Breiman (2001). In our analysis, variable importance is quantified as the reduction in $R^2_{oos}$ resulting from setting each variable, one at a time, to zero (its mean value post-normalization), with this metric normalized across all variables. For each method, the most significant variable, as per this definition, is assigned a value of one, and a color gradient is employed to visually represent the relative importance of each variable.

Table 1: Out-of-sample R-squared Values in Empirical Studies

<table>
<thead>
<tr>
<th></th>
<th>Ridge</th>
<th>Lasso</th>
<th>OLS/Ridgeless</th>
<th>RF</th>
<th>GBRT</th>
<th>NN($\ell_2$)</th>
<th>NN($\ell_1$)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Finance</td>
<td>0.80</td>
<td>-12.19</td>
<td>-81.08</td>
<td>-1.08</td>
<td>-14.21</td>
<td>0.30</td>
<td>-10.77</td>
</tr>
<tr>
<td>Finance</td>
<td>0.19</td>
<td>0.10</td>
<td>-1.25</td>
<td>0.08</td>
<td>-0.30</td>
<td>0.26</td>
<td>0.14</td>
</tr>
<tr>
<td>Macro 1a</td>
<td>15.29</td>
<td>15.40</td>
<td>-1375</td>
<td>24.39</td>
<td>16.44</td>
<td>16.94</td>
<td>19.09</td>
</tr>
<tr>
<td>Macro 1b</td>
<td>3.49</td>
<td>3.69</td>
<td>-2939</td>
<td>8.01</td>
<td>1.11</td>
<td>7.09</td>
<td>5.39</td>
</tr>
<tr>
<td>Macro 2</td>
<td>6.58</td>
<td>-14.58</td>
<td>-837</td>
<td>9.67</td>
<td>1.28</td>
<td>4.00</td>
<td>1.92</td>
</tr>
<tr>
<td>Micro 1</td>
<td>0.48</td>
<td>-1.01</td>
<td>-13198</td>
<td>-10.25</td>
<td>-5.07</td>
<td>0.49</td>
<td>-6.77</td>
</tr>
<tr>
<td></td>
<td>(0.84)</td>
<td>(2.01)</td>
<td>(12479)</td>
<td>(5.08)</td>
<td>(6.60)</td>
<td>(0.27)</td>
<td>(17.87)</td>
</tr>
<tr>
<td>Micro 2a</td>
<td>26.27</td>
<td>20.37</td>
<td>-12729</td>
<td>27.63</td>
<td>16.44</td>
<td>23.87</td>
<td>23.37</td>
</tr>
<tr>
<td></td>
<td>(7.50)</td>
<td>(6.41)</td>
<td>(9213)</td>
<td>(6.10)</td>
<td>(3.40)</td>
<td>(10.07)</td>
<td>(10.09)</td>
</tr>
<tr>
<td>Micro 2b</td>
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<td>-3.43</td>
<td>-14724</td>
<td>0.72</td>
<td>-6.45</td>
<td>3.16</td>
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<td>(5.25)</td>
<td>(10506)</td>
<td>(2.41)</td>
<td>(6.83)</td>
<td>(4.52)</td>
<td>(5.09)</td>
</tr>
</tbody>
</table>

Note: This table reports $R^2_{oos}$ values, presented in percentages, for Ridge, Lasso, OLS/Ridgeless, RF, GBRT, and NNs with respective $\ell_1$ and $\ell_2$ penalties, across six empirical studies spanning Finance, Macroeconomics, Microeconomics. For the first example in Macroeconomics and the second example in Microeconomics, two benchmark models are considered for comparison. Where standard deviations are applicable, they are provided in parentheses.
Figure 10: Variable Importance Plots

Finance 1

Finance 2

Macro 1b

Macro 2

Micro 1

Micro 2b

Note: This figure illustrates the variable importance across six empirical studies, using color gradients to show the relative reductions in $R^2_{oos}$ by each covariate. For the first example in Macroeconomics and the second example in Microeconomics, we only present the cases with a more complex benchmark model.
4.1 Finance 1: Market Equity Premium

In the first analysis, we focus on predicting market equity returns using a dataset of financial and macroeconomic indicators compiled by Welch and Goyal (2007). This dataset comprises 16 predictors and includes 74 annual observations, covering a period from 1948 to 2021. Despite Welch and Goyal (2007) reporting a consistently negative $R^2_{oos}$ for this dataset, several other studies, such as those by Campbell and Thompson (2007), Ferreira and Santa-Clara (2011), Rapach et al. (2010), Kelly and Pruitt (2013), and Kelly et al. (2023), have developed forecasting strategies resulting in economically meaningful $R^2_{oos}$ values. These modest $R^2_{oos}$s translate into significant economic gains through simple market timing strategies, as extensively discussed in these aforementioned studies.

We revisit this exercise. The procedure involves selecting an expanding window of in-sample data, using cross-validation for optimal tuning parameter selection, and then refitting the model for predictions on a test sample. Following the empirical framework of Giannone et al. (2022), the initial training set spans from 1948 to 1964, with the model’s performance evaluated using the 1965 data. Subsequently, data from 1965 is added to the training set, and the process is repeated, with the next model tested on the 1966 data. This procedure is conducted 57 times in total, progressively incorporating an additional year’s data into the training set and shifting the test sample forward by one year each time.

We evaluate $R^2_{oos}$ from 57 different predictions. Despite these predictions stemming from 57 distinct models, the empirical results, as detailed in Table 1, corroborate our theoretical predictions. Specifically, Ridge regression records an $R^2_{oos}$ of 0.80%, significantly outperforming Lasso’s -12.19%. With the smallest sample size being 17, which is adequate for running OLS with 16 predictors, the OLS model notably exhibits a highly negative $R^2_{oos}$ of -81.08%. In comparison, the NN with an $\ell_2$ penalty, NN($\ell_2$), attains an $R^2_{oos}$ of 0.30%, while the same model with an $\ell_1$ penalty, NN($\ell_1$), shows a lower $R^2_{oos}$ of -10.77%. As indicated in Figure 10, Ridge and NN($\ell_2$) appear to assign similar weights to these covariates. The leading covariate, eqis—the equity issuing activity ratio—is closely followed in importance by the dividend-price ratio, d/p. The comparatively modest $R^2_{oos}$ of NNs, especially when compared to Ridge, could be attributed to the complexity of the model, which elevates the risk of overfitting, especially since our NN comprises 72 parameters. The low signal-to-noise ratio in this scenario hinders both tree algorithms from achieving a positive $R^2_{oos}$. Neverthe-

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16 The data was sourced from Amit Goyal’s website, accessible at https://sites.google.com/view/agoyal145, and processed using the methodology provided by Giannone et al. (2022).

17 In this example, $R^2_{oos} = 1 - \frac{\sum_{t=1965}^{2021} (y_t - \bar{y}_t)^2}{\sum_{t=1965}^{2021} (y_t - \bar{y}_t)^2}$, where $\bar{y}_t = \sum_{s=1948}^{t-1} y_s / (t - 1948)$.
less, RF outperforms GBRT, with the latter registering the lowest $R^2_{oos}$ among all evaluated machine learning algorithms.

4.2 Finance 2: Cross-Section of Expected Returns

In our second analysis, we build upon the predictors utilized by Gu et al. (2020) for predicting equity returns, extending the data up to December 2021. We analyze monthly total individual equity returns from the CRSP database for all firms listed on the NYSE, AMEX, and NASDAQ. The average number of stocks analyzed per month exceeds 6,200. Our dataset starts from March 1957, and we compile a total of 920 covariates to predict future returns. Following Gu et al. (2020), our initial training phase utilizes data from 1957 to 1986, followed by performance evaluation using 1987 data. This process is repeated 35 times, with each iteration expanding the training sample by an additional year and shifting the evaluation period forward by one year.

Table 1 compares the model performance in terms of $R^2_{oos}$, where the zero estimator serves as the benchmark, according to the recommendation by Gu et al. (2020). In this case, NN($\ell_2$) emerges as the leading model, closely followed by Ridge, achieving $R^2_{oos}$ of 0.26% and 0.19%, respectively. These models dominate NN($\ell_1$) and Lasso, which record an $R^2_{oos}$ of 0.14% and 0.10%. The performance of the OLS continues to be underwhelming in this exercise. As for the tree-based models, RF demonstrates superior performance compared to GBRT. The former achieves a slightly positive $R^2_{oos}$, whereas the latter exhibits a slight negative $R^2_{oos}$. Despite their inherent ability to capture non-linear relationships and interactions, the effectiveness of tree models is somewhat limited in scenarios characterized by low signal strength. Figure 10 reveals an intriguing pattern: there appears to be a relationship between the relatively stronger performance among these pairs — Ridge vs Lasso, RF vs GBRT, and NN($\ell_2$) vs NN($\ell_1$) — and their respective patterns of sparsity in variable importance plots. Models with denser variable weights outperform those with sparser ones.

To illustrate the economic significance of these relatively low $R^2_{oos}$s, we adopt the approach outlined by Gu et al. (2020) and devise a stock selection portfolio strategy. This strategy involves going long on the top 10% and shorting the bottom 10% of stocks, sorted based on their predicted returns for the upcoming month, with equal weighting applied to each stock.

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18 Due to the large scale of the dataset and the resulting computational limitations, we have adopted a two-fold cross-validation approach for this analysis. Additionally, a more restrictive grid selection is employed for the Lasso and Ridge models. Specifically, for Ridge, log($\lambda$) is set to range between 6 and 7, while for lasso, it varies from -3.4 to -2.4. The optimal tuning parameters fall within the central range of these specified grids.

19 In this pooled regression setting, we define $R^2_{oos} = 1 - \sum_{t \in OOS} (y_{i,t} - \hat{y}_{i,t})^2 / \sum_{t \in OOS} y_{i,t}^2$. 

Electonic copy available at: https://ssrn.com/abstract=4722678
every month. In terms of performance, NN(ℓ₂) achieves the highest Sharpe ratio at 2.13, indicating its superior risk-adjusted returns. This is closely followed by Ridge regression with a Sharpe ratio of 1.64. NN(ℓ₁) also demonstrates commendable performance, yielding a Sharpe ratio of 1.55. GBRT exhibits the least impressive performance, with the lowest Sharpe ratio of 0.80, which aligns with its underwhelming predictive performance.

4.3 Macro 1: Macroeconomic Forecasting

The prediction of US macroeconomic activity using a wide range of predictors has been a topic of significant interest since its initial exploration by Stock and Watson (2002). In our current study, we utilize the FRED-MD dataset, compiled by McCracken and Ng (2016), to forecast the monthly growth rate of US industrial production (IP). This dataset includes 119 potential predictors, covering a diverse array of macroeconomic indicators, and extends from February 1960 to December 2019. Our evaluation methodology aligns with the prediction procedure outlined by Giannone et al. (2022). We begin by training these machine learning models using data from February 1960 to December 1974 and then evaluate their performance on data from the subsequent year. This process is repeated 45 times, with each iteration expanding the training dataset by one year (12 observations) and similarly shifting the evaluation period forward. We adhere to the guidelines set by McCracken and Ng (2016) for transforming the covariates. Additionally, we follow their prescribed approach for managing data quality issues, which involves the removal of outliers and the filling of missing data.

We initiate our analysis with a benchmark model that includes only an intercept term. In this scenario, all machine learning models significantly outperform this benchmark, achieving $R^2_{oos}$ values ranging from 14.13% to 24.15%. On the other hand, the OLS model, somewhat expectedly, overfits the data, resulting in a negative $R^2_{oos}$ of -16. This outcome suggests the presence of strong signal strength within a high-dimensional set of covariates. The benchmark model’s lack of competitiveness aligns with our expectations, particularly when considering the temporal dependence prevalent in macroeconomic time series. Therefore, a more suitable benchmark model should incorporate lagged values of IP growth.

We thereby propose an alternative benchmark that incorporates an Autoregressive (AR) component. Within each training sample, we fit an AR model to the IP growth, selecting its order based on the AIC. The residuals from this model then serve as our prediction target. As discussed in Section 2.8, this approach effectively combines the predictions from the AR model with those from our machine learning models, yielding a hybrid output out-of-

\footnote{In this case, the definition of $R^2_{oos}$ is similar to how it is defined in the Finance 1 case.}
sample. Consequently, the new benchmark for comparison becomes the direct use of the AR model’s predictions, where adding zero implies no alteration to the prediction. In this alternative setup, the comparison of $R^2_{oos}$ values reveals a pattern somewhat associated with scenarios of weak signal strength: NN($\ell_2$) and RF emerge as the top performers, achieving $R^2_{oos}$ values of 7.09% and 8.01%, respectively. Following closely are NN($\ell_1$) at 5.39%, while GBRT lags with a considerably lower $R^2_{oos}$ of 1.11%. This disparity in performance appears associated with the findings in Figure 10, which illustrates GBRT’s tendency towards sparser models in comparison to their counterparts. In this case, linear models, specifically Ridge and Lasso, demonstrate comparable performance, achieving $R^2_{oos}$ values of 3.49% and 3.69%, respectively. This pattern suggests that the primary challenges associated with signal weakness are most evident from nonlinear features.

### 4.4 Macro 2: Economic Growth Across Countries

Next, we explore a dataset originally compiled by Barro and Lee (1994), which includes 60 socio-economic, institutional, and geographical covariates across 90 countries. This dataset is utilized for predicting long-term economic growth, specifically measured by the growth rate of GDP per capita from 1960 to 1985. A pivotal aspect of this analysis involves testing a key prediction of the classical Solow-Swan-Ramsey growth model, which concerns the effect of an initial (lagged) GDP per capita level on subsequent growth rates. By incorporating the logarithm of each country’s GDP per capita in 1960 alongside a constant term, our prediction model includes a total of 62 potential covariates.

Belloni et al. (2013b) implement the Square-root-Lasso technique in their regression, anticipating sparsity among the control variables. This methodology results in a remarkable sparse model, characterized by the inclusion of a singular control variable: the log of the black market premium, a measure of trade openness. In contrast, Giannone et al. (2022) employ a Bayesian approach with a spike-and-slab prior, concluding that a dense model, which includes all covariates, yields the best log-predictive score. In our predictive analysis, we adopt the same empirical methodology outlined by Giannone et al. (2022). We begin by randomly selecting half of the data samples for model estimation and then proceed to assess the performance of these models using the remaining

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21In implementing advanced machine learning models with a hybrid benchmark linear component $W\gamma$, we adopt a methodology that parallels the one used in Eq. (14). This approach entails a DGP assumption that $M_{W'}y$ is a general function of $M_{W'}X$. This assumption plays a critical role in streamlining the implementation of these machine learning methods, ensuring that the results are directly comparable to those obtained in linear settings. However, it is important to note that this DGP assumption is generally not equivalent to the assumption that $y - W\gamma$ is a function of $X$. 

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samples. This process is repeated 100 times. The average out-of-sample $R^2_{oos}$ from these 100 repetitions, in comparison to a benchmark model that includes only the intercept, is presented in Table 1, accompanied by their standard deviations, provided in parentheses.\textsuperscript{22,23}

Our empirical findings align with those of Giannone et al. (2022), indicating a similar pattern that dense models, specifically Ridge, RF, and NN($\ell_2$), exhibit superior performance compared to their sparse counterparts, such as Lasso, GBRT, and NN($\ell_1$). The limited sample size appears to disadvantage complex NN models, rendering them less effective than the simpler Ridge regression. RF demonstrates strong performance, achieving an $R^2_{oos}$ of 9.67%, although it concurrently introduces a twofold increase in the variability of $R^2_{oos}$ values compared to those based on Ridge regression. The variance of Lasso is pronounced, driven by a handful of extreme values; excluding these anomalies, its $R^2_{oos}$ improves to -0.56%. Across all evaluated models, the black market premium consistently emerges as the most influential variable in Figure 10, aligning with the sole variable selected by Belloni et al. (2013b).

4.5 Micro 1: Crime Rates across US States

Our first microeconomic case revisits the study by Donohue and Levitt (2001), which analyzes the effect of the legalization of abortion following the Roe vs. Wade decision in 1973 on the decline in crime rates. Their dependent variable is the change in log per-capita murder rates from 1986 to 1997 across 48 states, with a total of 576 observations. This variable is then regressed on the effective abortion rate. To account for potential confounding factors, Belloni et al. (2013a) expanded the control set used by Donohue and Levitt (2001) by including interactions and higher-order terms, resulting in a comprehensive set of 284 variables.

When Belloni et al. (2013a) employ the Lasso method for the selection of control variables for murder rate in their analysis, they discover that none of the control variables were selected.\textsuperscript{24} In a similar vein, Giannone et al. (2022) observe from their Bayesian analysis of this regression that the posterior density is concentrated on very low probability values of the slab component, which suggests that the regression model is sparse with high likelihood.

\textsuperscript{22}Here and thereafter the definition of $R^2_{oos}$ is standard: $R^2_{oos} = 1 - \sum_{i \in OOS} (y_i - \hat{y}_i)^2 / \sum_{i \in OOS} (y_i - \bar{y})^2$, where $\bar{y}$ is based on in-sample average of $y$.

\textsuperscript{23}We may also consider a benchmark model with GDP per capita in 1960 included, as predicted by theory. Interestingly, enforcing the inclusion of this variable in the model leads to a reduction in predictive performance across all models. In essence, adding this variable leads to a negative $R^2_{oos}$ compared to the model that includes only an intercept.

\textsuperscript{24}Belloni et al. (2013a) proposes a double-Lasso estimator to make inference on the effect of abortion on murder rate. Part of their procedure involves a Lasso regression of murder rate on control variables. It is important to note that they use differences as the dependent variable, but observe no substantial changes when using levels instead.
In a recent study, Guo and Toulis (2023) employ a randomization test to assess the null hypothesis that all regression coefficients are zero. Their test fails to reject this hypothesis.

We employ the same benchmark model and sample splitting strategy outlined by Giannone et al. (2022). For the initial estimation, we use data spanning from 1986 to 1989, covering all states. Additionally, we incorporate data from a randomly selected 50% of the states for the period from 1990 to 1997. The remaining 50% of the states from 1990 to 1997 are set aside for evaluating the model. This procedure is iterated 8 times, with each iteration expanding the training sample to include one additional year of data, starting from 1990, while correspondingly adjusting the evaluation sample to reflect this change. The entire sequence is carried out 13 times in total, yielding $8 \times 13 = 104$ distinct training and evaluation samples. We report the mean and standard deviation of $R^2_{oos}$ in Table 1.

Our findings reveal that NN($\ell_2$) exhibits a slight edge over Ridge regression, attaining a marginally superior $R^2_{oos}$ of 0.49%, compared to Ridge’s 0.48%. Apart from these two models, all other models tested demonstrate negative $R^2_{oos}$ values, which suggests the presence of very weak signals in the data. Our argument posits that the scarcity of significant signals observed in the existing literature is likely attributable to signal weakness. The empirical evidence does not definitively categorize the underlying DGP as either dense or sparse. It may simply be that no individual signals are particularly strong. While in such scenarios, a sparse model could seem like a reasonable approximation, our results reveal that the cumulative predictive power of weak signals, though individually insubstantial, is collectively non-negligible. This is corroborated by Figure 10, where we observe that both Ridge and NN($\ell_2$) assign small weights to nearly all covariates. Although RF follows a similar pattern, our simulations indicate that it is more adversely affected by the weak signals compared to the other models.

4.6 Micro 2: Eminent Domain and Economic Outcomes

In our final study, we concentrate on a regression setting pertinent to eminent domain. Previous research by Chen and Yeh (2012), and subsequently Belloni et al. (2012), employ instrumental variable regressions to understand the impact of eminent domain on economic outcomes. Differing from their broader focus, our study aligns closely with Giannone et al. (2022), who concentrate on the first stage of this regression. This involves predicting pro-plaintiff decisions in takings law cases based on the characteristics of judicial panels. Their dataset includes 138 potential covariates and a total of 312 observations.

Adopting their strategy, we estimate the model using data spanning from 1979 to 1984 for all circuits. This is augmented with data from 1985 to 2004, selected randomly for 50%
of the circuits. We assess the model’s performance with 1985 data from the circuits not included in the training set. Since the period from 1985 to 2004 encompasses 20 years, we repeat this procedure 20 times. Each repetition involves a new random selection of half of the circuits and the sequential addition of one year’s data to the training set, while correspondingly updating the evaluation set. This entire process is independently executed five times, resulting in a total of $20 \times 5 = 100$ distinct training and evaluation datasets.

In our analysis, we initially consider a benchmark model that includes only an intercept. In this setting, all machine learning models successfully identify predictive signals, as evidenced by significant $R^2_{\text{OOS}}$. RF emerges as the top performer with an $R^2_{\text{OOS}}$ of 27.63%, with other models also showing strong results, albeit GBRT being an exception. However, the scenario shifts markedly when the benchmark model is expanded to include not only the intercept but also additional variables. These include a dummy variable for the absence of cases in a given circuit-year and the number of takings appellate decisions, bringing the total to three covariates. Against this simple benchmark, the incremental predictive power contributed by the remaining covariates diminishes dramatically. Ridge’s $R^2_{\text{OOS}}$ falls to 1.89%, that of RF to 0.47%, and NN($\ell_2$) to 3.16%, with the $R^2_{\text{OOS}}$ of all other methods turning negative. Intriguingly, as highlighted in Figure 10, these results seem to associate with the distinct approaches these methods take in weighting covariates. Ridge, RF, and NN($\ell_2$) assign small weights uniformly across all covariates. Meanwhile, NN($\ell_1$) also opts for a model with a considerable number of coefficients, resulting in a performance that slightly surpasses both Lasso and GBRT, which favor more sparse models in this case.

5 Conclusion

In this paper, we scrutinize the performance of machine learning techniques in contexts characterized by low signal-to-noise ratios, a situation frequently observed in economics and finance. Our theoretical analysis indicates that while Lasso is often considered a modern alternative to traditional ordinary least squares, its application in these areas should be approached cautiously, primarily due to its lessened effectiveness with weak signals.

Our research complements and expands upon the arguments made by Giannone et al. (2022), who cast doubt on the prevalence of sparsity in economic datasets. We take this debate further by showing that it is signal weakness, not necessarily the absence of sparsity, that more significantly contributes to the observed limitations of Lasso in economic applications. Furthermore, the lack of significant variables in empirical studies may be attributed
more to signal weakness than to the sparse nature of the underlying DGP.

Our analysis also reveals a marked difference in the performance of Ridge regression. Notably, Ridge demonstrates superior resilience and effectiveness in these environments. Our theoretical findings are further substantiated by simulation studies encompassing a range of advanced machine learning techniques, including trees and neural networks. These experiments consistently reveal that algorithms designed to exploit sparsity tend to underperform in environments where signals are inherently weak. Broadly, our findings emphasize the importance of a nuanced, context-sensitive application of machine learning techniques, adapting to the distinctive data characteristics encountered across various domains.

References


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Online Appendix of
Can Machines Learn Weak Signals?

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February 11, 2024

Abstract

Appendix A validates the robustness of our theoretical predictions through Monte Carlo simulations. Appendix B provides an in-depth discussion on the selection of tuning parameters. Appendix C contains mathematical proofs of the main theorems of the paper. Appendix D is dedicated to the exposition of technical lemmas and their corresponding proofs.

A Supplemental Simulation Results

In our subsequent series of experiments, we intentionally deviate from the assumptions originally established during the development of our theoretical framework. This deviation is aimed at evaluating the robustness and generalizability of our theoretical predictions beyond their premises and initial parameters. To facilitate this evaluation, we introduce specific modifications to the baseline configuration along three distinct dimensions: First, we raise the proportion of zeros from 50% in the benchmark scenario to 90% in the DGP of betas. This alteration creates a scenario more favorable to Lasso compared to the baseline. Second, we amplify the signal-to-noise ratio by reducing the $p/n$ ratio—accomplished by decreasing $p$ while maintaining $n$—thereby offering an edge to both Ridge and Lasso over the zero estimator. Third, we modify the distribution of $Z$ from standard Gaussian to a t-distribution.

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characterized by four degrees of freedom, and with a mean of zero and a variance of one. In addition, we introduce heteroscedasticity into the error distribution, following the configuration outlined by Giannone et al. (2022). The error term’s variance is defined by the function \( \sigma^2 \exp(\alpha X_i^\top \delta / \sqrt{\sum_{i=1}^{n} (X_i^\top \delta)^2 / n}) \) with \( \alpha = 0.5 \). Here, \( X_i \) represents the i-th row of \( X \). \( \sigma \) serves as a scaling parameter to standardize the variance and match \( \sigma^2_\varepsilon = 1 \). The vector \( \delta \) is a \( p \times 1 \) vector with zero elements in the same positions as the zero elements of \( \beta_0 \), while non-zero elements are drawn from a standard Gaussian distribution.

Figure 1 provides a comparison of histograms for the various cases under consideration. In Case I, the Lasso estimator exhibits improved performance compared to the preceding baseline scenario. This improvement is evident as a shift of probability mass occurs to the left side of zero. Such a shift suggests that, for these particular realizations, Lasso surpasses the zero estimator in predictive accuracy. However, the Ridge estimator consistently outperforms the zero estimator across nearly all realizations. Therefore, despite the presence of a high proportion (90%) of zero coefficients in this particular DGP, Lasso remains a less favorable choice when compared to both the zero and Ridge estimators. In Case II, we observe slight improvements in the performance of Ridge and Lasso in comparison to the zero estimator. This seemingly modest change is attributed to the scaling of the relative error by \( np^{-1} \tau^2 \). Absent this scaling factor, the relative error would manifest as double the negative magnitude relative to the baseline scenario. On the other hand, Case III demonstrates the robustness of our theoretical findings, as it aligns closely with the baseline scenario despite variations in distributional assumptions.

B Choice of Tuning Parameters

Table 1 provides details regarding model configuration and tuning parameters. For Ridge and Lasso methods, which each involve only one tuning parameter, we employ the glmnet package. This package effectively determines the optimal tuning parameter through a default ten-fold cross-validation process. The process is conducted on an adaptively selected grid, ensuring efficient and effective selection of the optimal tuning parameter. Regarding our implementation of RF, GBRT, and NNs, we follow the protocol outlined in the simulation section. In the case of NNs, we adhere to a uniform architectural choice across our analyses, featuring a single hidden layer. The number of neurons in this hidden layer is approximately equal to the square root of the total number of neurons in the input layer, aligning the architecture with the complexity and dimensions of the dataset. By not tuning the NN
Figure 1: Robustness Analysis of Ridge and Lasso in Alternative DGPs

Note: The histograms illustrate the relative prediction error $\Delta(\hat{\beta}(\lambda_n))$ for Ridge (upper panel) and Lasso (lower panel) based on 1,000 Monte Carlo samples. We explore three distinct DGPs, each involving the alteration of a specific condition. For the left panel, we set $q = 0.1$. In the middle panel, we adjust $n/p$ to 0.5. In the right panel, we introduce t-distributed covariates with heterogeneous variance of $\varepsilon$. The benchmark DGP adheres to the following specifications: $n = 2,000$, $q = 0.5$, $n/p = 1$, and complies with Assumptions 1 and 2. For Ridge, we employ a fixed value of $\lambda = 1$, while for Lasso, we select a constant $C_\lambda = 5$ throughout these experiments. In each of the figures, vertical red dashed lines serve the same interpretive purpose as those in Figures 5 and 7. The benchmark cases for Ridge and Lasso are represented by the plots in the middle of the second row of Figures 5 and 7.

architecture extensively, we streamline the model selection process while retaining adequate complexity for effective learning. For the remaining tuning parameters in trees and NNs, we select suitable ranges based on model performance from the cross-validation step. A critical element in selecting our grid is to ensure that the optimal tuning parameters are situated within the median range of the grid.

C Mathematical Proofs

C.1 Proof of Theorem 1

Proof. Throughout the proof, we employ the shorthand notation "w.a.p.1" to denote "with probability approaching one."

For convenience, we omit the subscript $F$ from the expectation operator $\mathbb{E}_F(\cdot)$. Our
Additionally, we fix the batch size for SGD and focus on jointly tuning the learning rate (lr) and the number of epochs (epochs), as well as the \( \ell_1 \)- or \( \ell_2 \)-penalty parameter (log(\( \lambda \))).

Note: The table reports the range of tuning parameters for RF, GBRT, and NNs, as well as the architecture of NNs applied across six datasets. For RF, we fix the number of trees at \#trees= 500, and tune two other parameters: the depth of the tree (depth) and the number of features (#features) within a predefined grid. In the case of GBRT, we tune not only depth and #trees, but also the learning rate (lr). For NNs, we adopt a fixed model architecture, denoted by the number of neurons in each layer indicated in brackets. Additionally, we fix the batch size for SGD and focus on jointly tuning the learning rate (lr) and the number of epochs (epochs), as well as the \( \ell_1 \)- or \( \ell_2 \)-penalty parameter (log(\( \lambda \))).

<table>
<thead>
<tr>
<th>Dataset</th>
<th>Depth</th>
<th>Trees</th>
<th>Features</th>
<th>Architecture</th>
<th>( \ell_2 ) Parameters</th>
<th>( \ell_1 ) Parameters</th>
</tr>
</thead>
<tbody>
<tr>
<td>Finance 1</td>
<td>1~6</td>
<td>500</td>
<td>1~10</td>
<td>architecture={119,8,1}</td>
<td>(lr,epochs)∈{(0.1,2), (0.05,250), (0.005,500)}</td>
<td>log(( \lambda )) ∈ [−2, 1]</td>
</tr>
<tr>
<td>Macro 1</td>
<td>5~40</td>
<td>500</td>
<td>2~40</td>
<td>architecture={119,8,1}</td>
<td>(lr,epochs)∈{(0.0005,10), (0.002,50), (0.005,200)}</td>
<td>log(( \lambda )) ∈ [−10, 1.5]</td>
</tr>
<tr>
<td>Macro 2</td>
<td>1~10</td>
<td>500</td>
<td>2~60</td>
<td>architecture={297,16,1}</td>
<td>(lr,epochs)∈{(0.001,10), (0.005,250), (0.01,1000)}</td>
<td>log(( \lambda )) ∈ [−7, 4]</td>
</tr>
<tr>
<td>Micro 1</td>
<td>1~5</td>
<td>500</td>
<td>1~5</td>
<td>architecture={217,16,1}</td>
<td>(lr,epochs)∈{(0.0008,10), (0.0001,100), (0.0005,2000)}</td>
<td>log(( \lambda )) ∈ [−11, −7]</td>
</tr>
<tr>
<td>Micro 2</td>
<td>1~6</td>
<td>500</td>
<td>1~30</td>
<td>architecture={217,16,1}</td>
<td>(lr,epochs)∈{(0.01,10), (0.005,200), (0.001,1000)}</td>
<td>log(( \lambda )) ∈ [−12, −9]</td>
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<tr>
<td>Micro 2b</td>
<td>1~5</td>
<td>500</td>
<td>1~5</td>
<td>architecture={215,16,1}</td>
<td>(lr,epochs)∈{(0.001,10), (0.005,200), (0.01,100)}</td>
<td>log(( \lambda )) ∈ [0, 2]</td>
</tr>
</tbody>
</table>
objective is to demonstrate that
\[
\frac{\mathbb{E}\|\Sigma_2^{1/2}(\mathbb{E}(\beta_0|X, y) - \beta_0)\|^2}{\mathbb{E}\|\Sigma_2^{1/2}\beta_0\|^2} \to 1, \text{ as } n \to \infty.
\]

This can be shown by proving \(\mathbb{E}\|\Sigma_2^{1/2}\mathbb{E}(\beta_0|X, y)\|^2 = o(\mathbb{E}\|\Sigma_2^{1/2}\beta_0\|^2)\). Given that the eigenvalues of \(\Sigma_2\) are bounded away from zero and positive infinity, it suffices to establish that \(\mathbb{E}\|\beta_0|X, y\|^2 = o(\tau)\). Therefore, we need to prove for all \(1 \leq i \leq p\), \(\mathbb{E}(\beta_{0,i}|X, y)^2 = o(p^{-1}\tau)\), or, equivalently, \(\mathbb{E}(\beta_{0,i}|X, y)^2 = o(1)\).

By the inequality \(\mathbb{E}(\mathbb{E}(A|\mathcal{F})) \leq \mathbb{E}(\mathbb{E}(A|\mathcal{G}))\) for \(\mathcal{F} \subset \mathcal{G}\), and that \(\beta_0\) is i.i.d., we have
\[
\mathbb{E}(\mathbb{E}(b_{0,i}|X, y)) \leq \mathbb{E}(\mathbb{E}(b_{0,i}|X, y, \beta_{0,-i})) = \mathbb{E}(\mathbb{E}(b_{0,i}|X, \beta_{0,i}, \Sigma^{-1/2}X, i + z))^2,
\]
where \(z\) is defined in Assumption 2, \(X_{:,i}\) represents the \(i\)-th column of \(X\), and \(\beta_{0,-i}\) denotes the subvector of \(\beta\) without the \(i\)th entry. Denote the information set generated by \(\{X_{:,i}, \beta_{0,i}, \Sigma^{-1/2}X, i + z\}\) as \(\mathcal{G}_i\). Additionally, for any \(M_1 < 0\), find \(M_2\) (a function of \(M_1\)) such that \(\mathbb{E}b_{0,i}(M_1 \leq b_{0,i} \leq M_2) = 0\). This is always feasible because \(\mathbb{E}b_{0,i} = 0\). By Cauchy-Schwarz inequality, we have
\[
\mathbb{E}(\mathbb{E}(b_{0,i}|\mathcal{G}_i))^2 \leq 3\mathbb{E}\left(\mathbb{E}(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)|\mathcal{G}_i)\right)^2 + 3\mathbb{E}\left(\mathbb{E}(b_{0,i}I(b_{0,i} > M_2)|\mathcal{G}_i)\right)^2.
\]

Now we prove for any given \(M_1\), \(\lim_{n \to \infty} \mathbb{E}(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)|\mathcal{G}_i)^2 = 0\). Since \(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)\) is bounded, so is \((\mathbb{E}(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)|\mathcal{G}_i)^2\). By the dominated convergence theorem, it suffices to show: \(\mathbb{E}(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)|\mathcal{G}_i) = o_p(1)\).

Write \(\tilde{x}_k = (\Sigma^{-1/2}X_{:,i})_k\) and \(\tilde{y}_k = \beta_{0,i}x_k + z_k\) for \(k = 1, \ldots, n\). By definition, we have
\[
\mathbb{E}(b_{0,i}I(M_1 \leq b_{0,i} \leq M_2)|\mathcal{G}_i) = \frac{\int bI(M_1 \leq b \leq M_2) \exp \left( -\frac{\sum_{k=1}^{n} (\tilde{y}_k - p^{-1/2}\tau^{1/2}\tilde{x}_k b)^2}{2} \right) dF(b)}{\int \exp \left( -\frac{\sum_{k=1}^{n} (\tilde{y}_k - p^{-1/2}\tau^{1/2}\tilde{x}_k b)^2}{2} \right) dF(b)} = \frac{\int bI(M_1 \leq b \leq M_2) \exp \left( -\frac{p^{-1}\tau b^2 \sum_{k=1}^{n} \tilde{x}_k^2 - 2bp^{-1/2}\tau^{1/2}\sum_{k=1}^{n} \tilde{y}_k \tilde{x}_k}{2} \right) dF(b)}{\int \exp \left( -\frac{p^{-1}\tau b^2 \sum_{k=1}^{n} \tilde{x}_k^2 - 2bp^{-1/2}\tau^{1/2}\sum_{k=1}^{n} \tilde{y}_k \tilde{x}_k}{2} \right) dF(b)}.
\]
where $F$ is the distribution function of $b_{0,i}$. Given that $b_{0,i}$ is a sub-exponential random variable, there exists a constant $C$ such that $\mathbb{E}\exp(C|b_{0,i}|)$ is finite. Define the event

$$A_n := \left\{ \left| p^{-1/2} \tau^{1/2} \sum_{k=1}^{n} \bar{y}_k \bar{x}_k \right| \leq C \right\},$$

and let

$$X_n(\bar{b}) := \bar{b} \mathbb{I}(M_1 \leq \bar{b} \leq M_2) \exp \left( -p^{-1} \tau \bar{b}^2 \sum_{k=1}^{n} \bar{x}_k^2 - 2\bar{b} p^{-1/2} \tau^{1/2} \sum_{k=1}^{n} \bar{y}_k \bar{x}_k \right),$$

where $\bar{b} \sim F$ is independent of $\bar{x}_k$ and $\bar{y}_k$, for $k = 1, \ldots, n$. It follows that

$$X_n(\bar{b}) 1_{A_n} - \bar{b} \mathbb{I}(M_1 \leq \bar{b} \leq M_2) 1_{A_n} \xrightarrow{P} 0, \quad \text{and} \quad \mathbb{P}(A_n) \to 1,$$

since by Lemmas 2 and 3, $p^{-1/2} \tau \bar{b}^2 \sum_{k=1}^{n} \bar{x}_k^2 = o_P(1)$ and $p^{-1/2} \tau^{1/2} \sum_{k=1}^{n} \bar{y}_k \bar{x}_k = o_P(1)$. Furthermore, since $|X_n(\bar{b}) 1_{A_n} - \bar{b} \mathbb{I}(M_1 \leq \bar{b} \leq M_2) 1_{A_n}| \leq |\bar{b} \exp(C|\bar{b}|)| + |\bar{b}|$, and the latter is integrable with respect to $F$, the dominated convergence theorem implies

$$\lim_{n \to \infty} \mathbb{E}|X_n(\bar{b}) 1_{A_n} - \bar{b} \mathbb{I}(M_1 \leq \bar{b} \leq M_2) 1_{A_n}| = 0,$$

which, along with the fact that $1_{A_n} \xrightarrow{P} 1$, lead to

$$\mathbb{E} \left[ |X_n(\bar{b}) - \bar{b} \mathbb{I}(M_1 \leq \bar{b} \leq M_2)||\bar{y}_k, \bar{x}_k, 1 \leq k \leq n \right] = o_P(1).$$

This further implies

$$\int \bar{b} \mathbb{I}(M_1 \leq b \leq M_2) \exp \left( -p^{-1} \tau \bar{b}^2 \sum_{k=1}^{n} \bar{x}_k^2 - 2\bar{b} p^{-1/2} \tau^{1/2} \sum_{k=1}^{n} \bar{y}_k \bar{x}_k \right) dF(b) \xrightarrow{P} \int \bar{b} \mathbb{I}(M_1 \leq b \leq M_2)dF(b) = 0.$$

In a similar vein, we can show that

$$\int \exp \left( -p^{-1} \tau \bar{b}^2 \sum_{k=1}^{n} \bar{x}_k^2 - 2\bar{b} p^{-1/2} \tau^{1/2} \sum_{k=1}^{n} \bar{y}_k \bar{x}_k \right) dF(b) \xrightarrow{P} \int dF(b) = 1.$$

As a result, (C1) implies

$$\lim_{n \to \infty} \mathbb{E}(\mathbb{E}(b_{0,i}|G_i))^2 \leq \lim_{n \to \infty} \mathbb{E}(\mathbb{E}(b_{0,i}\mathbb{I}(b_{0,i} > M_2)|G_i))^2 + 3 \lim_{n \to \infty} \mathbb{E}(\mathbb{E}(b_{0,i}\mathbb{I}(b_{0,i} < M_1)|G_i))^2$$

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This is because Eq. (C2) implies that
\[ \leq 3\mathbb{E}b_{0,i}^2 \mathbb{I}(b_{0,i} > M_2) + 3\mathbb{E}b_{0,i}^2 \mathbb{I}(b_{0,i} < M_1). \]

Since \( b_{0,i} \) has finite variance, the right hand side of the above inequality can be arbitrary small by letting \( M_1 \to -\infty \), which completes the proof. \( \square \)

### C.2 Proof of Theorem 2

**Proof.** For ease of notation, we let \( \hat{\beta} := \hat{\beta}_r(\lambda_n) \) and \( c_n := p/n \). Additionally, define \( \delta_1^* := 2\sqrt{\sigma_x^2 / \sigma_1^2}, \delta_2^* := (2\lambda \sigma_x^2 / \sigma_3^2 \theta_4 - 4\sigma_x^2 / \sigma_3^2 \delta_1^*) / \delta_1^1 \lambda, \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2^*) := (\delta_1^* \delta_2^* - 2\sigma_x^2 \sigma_3^2 \theta_4) / 4\sigma_3^2 \theta_3, \) and

\[
C_n^\phi := c_n \tau^{-1} \sigma_x^2 \sigma_\beta^2 - c_n \frac{\tau^{-1} \sigma_x^2 (\delta_1^1)^2}{4\lambda} + c_n \sigma_x^2 \sigma_3^1 \theta_3 - c_n \lambda \sigma_x^2 \sigma_3^2 \theta_4.
\]

We first show that it is sufficient to establish that
\[
c_n \tau^{-3/2}(\|\Sigma_{2}^{1/2}(\hat{\beta} - \beta_0)\| - \|\Sigma_{2}^{1/2} \beta_0\|) \xrightarrow{P} \alpha_2^* := \theta_2 \sigma_x^2 \left( \frac{\sigma_x^2 \theta_1}{2\lambda^2 \sigma_\beta} - \frac{\sigma_\beta}{\lambda} \right). \tag{C2}
\]

This is because Eq. (C2) implies that
\[
\|\Sigma_{2}^{1/2}(\hat{\beta} - \beta_0)\|^2 = \|\Sigma_{2}^{1/2} \beta_0\|^2 + 2c_n^{-1} \tau^{3/2} \alpha_2^* \|\Sigma_{2}^{1/2} \beta_0\| + o_p(c_n^{-1} \tau^2).
\]

On the other hand, using Lemma 2, we deduce that
\[
\|\Sigma_{2}^{1/2} \beta_0\| = \tau^{1/2} \sigma_x \sigma_\beta + O_p(p^{-1/2} \tau^{1/2}). \tag{C3}
\]

The above two equations together yield the result of the theorem.

To prove Eq. (C2), by incorporating (C3), it reduces to proving that
\[
c_n \tau^{-1}(\tau^{-1/2}\|\Sigma_{2}^{1/2}(\hat{\beta} - \beta_0)\| - \sigma_x \sigma_\beta) \xrightarrow{P} \alpha_2^*. \tag{C4}
\]

Set \( w = \tau^{-3/2} \Sigma_{2}^{1/2}(\hat{\beta} - \beta_0) \) and \( \hat{w} = \tau^{-3/2} \Sigma_{2}^{1/2}(\hat{\beta} - \beta_0) \). After rewriting Ridge’s optimization problem (2), \( \hat{w} \) equals
\[
\arg \min_w \frac{c_n}{n} \left( \tau^{-1/2}\Sigma_{2}^{1/2}Z w - \tau^{-1} \varepsilon \right)^2 + \sigma_x^2 \lambda \left( \Sigma_{2}^{1/2} w + \tau^{-3/2} \beta_0 \right)^2 - \frac{c_n \tau^{-2}}{n} \|\varepsilon\|^2 - C_n^\phi, \tag{C5}
\]

where subtracting \( c_n \tau^{-2} \|\varepsilon\|^2 / n \) and \( C_n^\phi \) from the objective function does not alter the solution. Using the definition of \( \hat{w} \), proving (C4) is equivalent to proving
\[ c_n \| \hat{w} \| - c_n \tau^{-1} n \sigma_x \sigma_\beta \xrightarrow{P} \alpha_2^*. \]  
\text{(C6)}

Equivalently, we need to prove for all \( \epsilon > 0 \), w.p.a.1,

\[ \alpha_2^* - \epsilon \leq c_n \| \hat{w} \| - c_n \tau^{-1} n \sigma_x \sigma_\beta \leq \alpha_2^* + \epsilon. \]  
\text{(C7)}

Next, we note from Lemma 13 that it suffices to prove the above convergence holds true for \( \hat{w}_B \), where \( \hat{w}_B \) is a solution to

\[
\arg \min_{w \in S_n^u} \frac{c_n}{n} \left( \tau^{1/2} \Sigma_1^{1/2} Z w - \tau^{-1} \| \varepsilon \|_2 \right)^2 + c_n^2 \lambda \left( \Sigma_2^{-1/2} w + \tau^{-3/2} \beta_0 \right)^2 - \frac{c_n}{n} \| \varepsilon \|_2^2 - C_n^\phi, \]  
\text{(C8)}

and \( S_n^u = \{ w | c_n \tau^{-1} n \sigma_x \sigma_\beta - K_\alpha \leq c_n \| w \| \leq c_n \tau^{-1} n \sigma_x \sigma_\beta + K_\alpha \} \) for some sufficiently large \( K_\alpha \).

We’ll denote the optimal solution as \( \hat{w} \) instead of using \( \hat{w}_B \) for simplicity.

Note that for any vector \( x \), \( \| x \|_2^2 = \max_{n} \sqrt{n} u \top x - n \| u \|_2^2 / 4 \), where its argmax is \( 2x / \sqrt{n} \), and similarly \( \| x \|_2^2 = \max_{n} v \top x - \| v \|_2^2 / 4 \). Applying these equalities to \( \| \tau^{1/2} \Sigma_1^{1/2} Z w - \tau^{-1} \| \varepsilon \|_2^2 \) and \( \| \Sigma_2^{-1/2} w + \tau^{-3/2} \beta_0 \|_2^2 \), setting \( \tilde{u} = \Sigma_1^{1/2} u \), and \( \tilde{v} = \Sigma_2^{-1/2} v \), we can rewrite (C8) as

\[
\min_{u \in S_w^u} \max_{\tilde{u}, \tilde{v}} \frac{c_n}{n} \tau^{1/2} \tilde{u} \top Z w - \frac{c_n}{n} \tau^{-1} \tilde{u} \top \Sigma_1^{1/2} \varepsilon - \frac{c_n}{n} \| \Sigma_1^{1/2} \tilde{u} \|_2^2 - \frac{2}{n} \| \varepsilon \|_2^2 - \frac{c_n}{n} \| \varepsilon \|_2^2 - C_n^\phi. \]  
\text{(C9)}

To simplify notation and without ambiguity, we continue using \( u \) and \( v \) in place of \( \tilde{u} \) and \( \tilde{v} \).

For a given \( w \), the argmax of Eq. (C9), denoted by \( \hat{u} \), is equal to \( \frac{1}{\sqrt{n}} (\tau^{1/2} Z w - \tau^{-1} \Sigma_1^{1/2} \varepsilon) \). Given the definition of \( S_w^u \) and Assumptions 1 and 2, we have \( \| w \| \leq \tau^{-1} n \sigma_x \sigma_\beta + c_n^{-1} K_\alpha \), \( \| \Sigma_1 \| \leq C_1 \), \( \| \varepsilon \| \leq C_\varepsilon \). Furthermore, w.p.a. 1, \( \| z \| \leq \sqrt{2n} \) by the law of large numbers, which implies \( \| \varepsilon \| \leq \sqrt{2C_\varepsilon n} \). Together with Lemma 6 and that \( \tau c_n \to 0 \) by Assumption 4, we have the following bound for \( \| \hat{u} \| \) as \( n \) is large enough:

\[ \| \hat{u} \| \leq \frac{2 \tau^{1/2}}{\sqrt{n}} \| \Sigma_1 Z w \| + \frac{2}{\sqrt{n}} \| \tau^{-1} \Sigma_1^{1/2} \varepsilon \| \leq 4 \tau^{-1} \sqrt{C_1 C_\varepsilon}. \]

Let \( S_u^v = \{ u | \| u \| \leq 4 \tau^{-1} \sqrt{C_1 C_\varepsilon} \} \). Based on the above result, w.p.a.1, the following optimization problem is equivalent to (C9):

\[
\min_{u \in S_w^u} \max_{v \in S_u^v} \frac{c_n}{n} \tau^{1/2} u \top Z w - \frac{c_n}{n} \tau^{-1} u \top \Sigma_1^{1/2} \varepsilon - \frac{c_n}{n} \| \Sigma_1^{1/2} u \|_2^2 - \frac{2}{n} \| \varepsilon \|_2^2 - \frac{c_n}{n} \| \varepsilon \|_2^2 - C_n^\phi. \]

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Next, we need introduce an auxiliary problem for the purpose of applying CGMT:

\[
\phi(g, h) = \max_{v \leq 4n^{-1} \sqrt{\Sigma}} \min_{w \in S_w} \max_{\|u\| = \delta} \mathcal{R}_n(w, v, u), \quad \text{where}
\]

\[
\mathcal{R}_n(w, v, u) = \frac{c_n \tau^{1/2}}{\sqrt{n}} \|w\| (g^\top u - \frac{c_n \tau - 1}{\sqrt{n}} \delta h^\top w - \frac{c_n \tau - 1}{\sqrt{n}} u^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n}{n} \|\Sigma_1^{-1/2} u\|^2) + \frac{c_n^2}{n} \|\Sigma_2^{1/2} v\|^2 - \frac{c_n^2 \lambda}{4} \varepsilon^2 - C_n^\phi,
\]

(C10)

and \( g \in \mathbb{R}^n \) and \( h \in \mathbb{R}^p \) are standard Gaussian vectors, independent of the other random variables. Similarly, let \( S_n := \{w \mid c_n \|w\| - c_n \tau - 1 \Sigma \sigma + \sigma \Sigma \} < \varepsilon \}, define \( \phi_{S_n}(g, h) \) as the optimal value of an analogous optimization problem to (C11), with \( w \) restricted to \( S_n \) \( \bar{\mathcal{S}}_n \).

Lemma 14 characterizes the limiting behavior of the optimal solution to (C10), \( \hat{w} \), and in turn, proves the desired (C6), under conditions pertaining to the optimization problem (C11). Therefore, we only need show that conditions outlined in Lemma 14 hold. That is, we need to prove the existence of the constants \( \bar{\phi} < \bar{\phi}_{S_n} \) such that for all \( \eta > 0 \), w.p.a.1 in the limit of \( n \to \infty \), \( \phi(g, h) < \bar{\phi} + \eta \) and \( \phi_{S_n}(g, h) > \bar{\phi}_{S_n} - \eta \).

Let \( \bar{u} = u/\delta \), maximizing part of \( \mathcal{R}_n(w, v, u) \) pertaining to \( u \) over \( u \) simplifies to the following problem:

\[
\max_{\|u\| = \delta} \frac{c_n \tau^{1/2}}{\sqrt{n}} \|w\| (g^\top u - \frac{c_n \tau - 1}{\sqrt{n}} u^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n}{n} \|\Sigma_1^{-1/2} u\|^2) = \max_{\|u\| = \delta} \frac{c_n \delta}{\sqrt{n}} (\tau^{1/2} \|w\| (g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon) \bar{u} - \frac{c_n \delta^2}{4} \bar{u}^\top \Sigma_1^{-1} \bar{u}).
\]

The latter is a quadratic programming problem, which has been extensively studied in, e.g., Gander et al. (1989) and Tao and An (1998). The optimal value associated with this problem is given by the following expression:

\[
-\frac{c_n^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon)^\top (\Sigma_1^{-1} - \mu_n(\alpha, \delta))^{-1} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon) (C12)
\]

where \( \alpha := \|w\| \) and \( \mu_n(\alpha, \delta) \) is the solution to

\[
\frac{1}{n} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon)^\top (\Sigma_1^{-1} - \mu_n(\alpha, \delta))^{-2} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon) - \frac{\delta^2}{4} = 0, \quad (C13)
\]
under the condition that \( \Sigma^{-1}_1 - \mu_n(\alpha, \delta)\mathbb{I} \) is positive semidefinite. Using this, Eq. (C11) can be rewritten as the following:

\[
\max_{0 \leq \delta \leq 4^{-1/4}\sqrt{c_1 c_2}} \min_{w \in S^w_0} -c_n \delta^2/4 \mu_n(\alpha, \delta) + \frac{c_n}{n} (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) \top (\Sigma^{-1}_1 - \mu_n(\alpha, \delta)\mathbb{I})^{-1} \\
\times (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) - \frac{c_n}{\sqrt{n}} \delta h \top w + c_n^2 \lambda v \top w \\
+ c_n^2 \lambda \tau^{-3/2} v \top \Sigma^{-1/2}_2 \beta_0 - \frac{c_n^2}{n} \lambda \|\Sigma^{-1/2}_2 v\|^2 \\
- \frac{c_n}{n} \|\varepsilon\|^2 - C_n^\phi.
\]

Solving the inside minimization problem with respect to \( w/\alpha \) while fixing \( \alpha \) leads to

\[
\max_{0 \leq \delta \leq 4^{-1/4}\sqrt{c_1 c_2}} \min_{v \in \mathbb{C}} \left| c_n \alpha - c_n \alpha - \tau^{-1} \gamma \sigma_\alpha \beta_0 \right| \leq K_\alpha \left| -c_n \delta^2/4 \mu_n(\alpha, \delta) + \frac{c_n}{n} (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) \top (\Sigma^{-1}_1 - \mu_n(\alpha, \delta)\mathbb{I})^{-1} \\
\times (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) - \frac{c_n}{\sqrt{n}} \delta h \top \alpha \\
+ c_n^2 \lambda \tau^{-3/2} v \top \Sigma^{-1/2}_2 \beta_0 - \frac{c_n^2}{n} \lambda \|\Sigma^{-1/2}_2 v\|^2 \\
- \frac{c_n}{n} \|\varepsilon\|^2 \right| + C_n^\phi.
\]

By Lemma 15, the objective function of the above optimization is convex in \( \alpha \) and jointly concave in \( (\delta, v) \). As a result, we can switch the order of min and max by Corollary 3.3 in Sion (1958). Also, note that for any vector \( x, \|x\| = \min_{\gamma > 0} \frac{1}{2\gamma} \|x\|^2 + \frac{\gamma}{2} \). Applying this equation to \( \|c_n \lambda v - n^{-1/2} \tau^{1/2} \delta h\| \alpha \), Eq. (C14) becomes

\[
c_n \alpha - \tau^{-1} \gamma \sigma_\alpha \beta_0 \leq K_\alpha \left| \max_{0 \leq \delta \leq 4^{-1/4}\sqrt{c_1 c_2}} \min_{\gamma > 0} -c_n \delta^2/4 \mu_n(\alpha, \delta) + \frac{c_n}{n} (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) \top \\
\times (\tau^{1/2}\alpha g - \tau^{-1}\Sigma^{-1/2}_1 \varepsilon) - \frac{c_n}{\sqrt{n}} \delta h \top \alpha \\
+ c_n^2 \lambda \tau^{-3/2} v \top \Sigma^{-1/2}_2 \beta_0 - \frac{c_n^2}{n} \lambda \|\Sigma^{-1/2}_2 v\|^2 \\
- \frac{c_n}{n} \|\varepsilon\|^2 \right| + C_n^\phi.
\]

Note that the objective function above is jointly concave in \( (\delta, \gamma, v) \). To see why this is true, it is sufficient to prove that \(-\frac{\alpha^2}{2\gamma}\|c_n \lambda v - n^{-1/2} \tau^{1/2} \delta h\|^2 \) is jointly concave in \( (\delta, \gamma, v) \), which follows by Lemma 13 in Thrampoulidis et al. (2018). Consequently, after solving the first maximization problem over \( v \), the resulting function remains jointly concave in \( (\delta, \gamma) \).

Maximizing over \( v \) is again a standard quadratic programming problem, which leads to
C.3 Proof of Theorem 3

Proof. For convenience, we define the shorthand notation \( \hat{\beta}_{i}^{j}(\lambda_{n}) \). Also, we define

\[
\hat{R}^{K-CV}(\lambda_{n}) := \frac{1}{n} \sum_{i=1}^{K} \| y(i) - X(i)\hat{\beta}_{i}^{j}(\lambda_{n}) \|^{2}. \tag{C17}
\]

By Lemma 6, w.p.a.1, we have
\[
\frac{1}{n} \| X_{(-i)}^T X_{(-i)} \| \leq \frac{1}{n} C_2 \| Z_{(-i)}^T Z_{(-i)} \| \leq C_2 (1 + \sqrt{c_n})^2, \quad i = 1, \ldots, K.
\] (C18)

Additionally, by Lemmas 2 and 3, w.p.a.1, we have
\[
\frac{1}{n} \| \varepsilon \|^2 \leq 2\sigma_\varepsilon^2 \quad \text{and} \quad \frac{1}{n} \| y \|^2 \leq 2\sigma_\varepsilon^2.
\] (C19)

Under the condition that \( \lambda_n \geq \epsilon \), using (C18) and (C19), we have, w.p.a.1,
\[
\| \hat{\beta}_{\lambda_n} - \hat{\beta}_{\lambda_n} \|^2 = \left\| \frac{1}{n} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_n \right)^{-1} X_{(-i)} y_{(-i)} \right\|^2 \leq \frac{\| X_{(-i)} y_{(-i)} \|^2}{c_n^2 \epsilon^2 n^2} \leq \frac{2 C_2 \sigma_\varepsilon^2 (1 + \sqrt{c_n})^2}{c_n^2 \epsilon^2}.
\] Using a similar argument, we have, w.p.a.1,
\[
\| \hat{\beta}_{\lambda_1} - \hat{\beta}_{\lambda_2} \|^2 = c_n^2 (\lambda_1 - \lambda_2)^2 \left\| \frac{1}{n} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_2 \right)^{-1} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_1 \right)^{-1} X_{(-i)} y_{(-i)} \right\|^2 \leq \frac{2 C_2 \sigma_\varepsilon^2 (1 + \sqrt{c_n})^2 (\lambda_1 - \lambda_2)^2}{c_n^2 \epsilon^2}.
\] (C20)

With the inequalities above and triangle inequalities, we obtain, w.a.p.1,
\[
\left| \hat{R}^{K-CV}(\lambda_1) - \hat{R}^{K-CV}(\lambda_2) \right| = \frac{1}{n} \left\| \sum_{i=1}^K \left( \| y(i) - X(i) \hat{\beta}_{\lambda_1} \|^2 - \| y(i) - X(i) \hat{\beta}_{\lambda_2} \|^2 \right) \right\| \leq \frac{2}{n} \sum_{i=1}^K \| y(i) \| \| \hat{\beta}_{\lambda_1} - \hat{\beta}_{\lambda_2} \| \left( \| y(i) \| + \frac{2 C_2 \sigma_\varepsilon (1 + \sqrt{c_n})}{c_n \epsilon} \| X(i) \| \right) \leq \tilde{C} | \lambda_1 - \lambda_2 |,
\] (C21)

where \( \tilde{C} \) is some fixed constant. Based on this inequality, Lemma 17 proves
\[
\inf_{\lambda \in [\epsilon, \tilde{c} \tau^{-1}]} \left\{ p n^{-1} \tau^{-2} \left( \hat{R}^{K-CV}(\lambda) - \frac{1}{n} \| \varepsilon \|^2 \right) - p n^{-1} \tau^{-2} \| \Sigma_{1/2}^{1/2} \beta_0 \|^2 \right\} > 0,
\] (C22)

w.p.a.1 for some constant \( \tilde{c} > 0 \). Additionally, as \( n \to \infty \), for any fixed \( \lambda > 0 \),
\[
p n^{-1} \tau^{-2} \left\{ \hat{R}^{K-CV}(\tau^{-1} \lambda) - \frac{1}{n} \| \varepsilon \|^2 - \| \Sigma_{1/2}^{1/2} \beta_0 \|^2 \right\} \to \frac{2(K-1)}{K} \theta_2 \sigma_x^4 \left( \frac{\sigma_\varepsilon^2}{2\lambda^2} - \frac{\sigma_\varepsilon^2}{\lambda} \right).
\] (C23)

Using (C23) and the definition of \( \lambda^{opt} \), we obtain, w.p.a.1,
\[
p n^{-1} \tau^{-2} \left\{ \hat{R}^{K-CV}(\tau^{-1} \lambda^{opt}) - \frac{1}{n} \| \varepsilon \|^2 - \| \Sigma_{1/2}^{1/2} \beta_0 \|^2 \right\}
\]
< 0 < \inf_{\lambda \in [\epsilon, \epsilon \tau^{-1}]} pn^{-1}\tau^{-2} \left\{ \hat{R}^{K-CV}(\lambda) - \frac{1}{n} \|\varepsilon\|^2 - \|\Sigma_2^{1/2} \beta_0\|^2 \right\}.

This suggests that the minimizer of \( \hat{R}^{K-CV}(\lambda) \) must satisfy \( \hat{\lambda}^{K-CV} \geq \tilde{c}\tau^{-1} \), w.a.p.1, so that, \( \hat{\lambda}^{K-CV}_n = \arg \min_{\lambda_n \in [\epsilon \tau^{-1}, \infty)} \hat{R}^{K-CV}(\lambda_n) \). Moreover, it also implies that \( \tau^{-1} \lambda^{opt} \notin [\epsilon, \tilde{c}\tau^{-1}] \), that is, \( \lambda^{opt} \geq \tilde{c} \). Next, we re-parametrize the above optimization problem:

\[
\tilde{\mu} = \arg \min_{\mu \in [0, \tilde{\epsilon}]} \tilde{R}(\mu), \quad \text{where} \quad \tilde{R}(\mu) := \hat{R}^{K-CV}(\tau^{-1} \mu^{-1}),
\]

and we extend the domain of \( \tilde{R}(\cdot) \) to include 0: \( \tilde{R}(0) := \lim_{\mu \rightarrow 0} \tilde{R}(\mu) = \|y\|^2/n \). Lemma 18 implies that \( pn^{-1}\tau^{-2}\tilde{R}(\mu) \) satisfies stochastic equicontinuity. Using this fact and Theorem 1 of Newey (1991), the convergence of

\[
lp^{-1}\tau^{-2} \left( \tilde{R}(\mu) - \frac{1}{n} \|\varepsilon\|^2 \right) - pn^{-1}\|\Sigma_2^{1/2} \beta_0\|^2 \overset{P}{\rightarrow} 2(K-1)K\theta_2 \sigma_x^4 \left( \sigma_x^2 \mu^2 - \sigma^2_\beta \mu \right)
\]

holds uniformly over the interval \([0, \tilde{c}\tau^{-1}]\). Since \( (\lambda^{opt})^{-1} \) is a unique minimizer of the right-hand-side and is distinct from zero, it follows that \( \tilde{\mu} \overset{P}{\rightarrow} (\lambda^{opt})^{-1} \) and \( \tilde{\mu} = \tau^{-1}(\hat{\lambda}^{K-CV}_n)^{-1} \), w.a.p.1, which conclude the proof.

### C.4 Proof of Theorem 4

**Proof.** For ease of notation, we let \( \hat{\beta} := \hat{\beta}_i(\lambda_n) \). We adopt the same notation \( \delta_1^* \) and \( \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \) as used in the proof of Theorem 2. With respect to \( \delta_2^* \) and \( C^\phi_n \), we define them as \( 2\sigma_x^2 \sigma_\beta^2 \theta_4/\delta_1^* \) and \( c_n \tau^{-1} \sigma_x^2 \sigma_\beta^2 \), respectively.

Analogous to the proof of Theorem 2, it is essential to establish that the following inequality holds w.a.p.1 for any sufficiently small \( \epsilon > 0 \):

\[
\frac{c_\alpha}{2\sigma_\beta} + \epsilon \leq c_n \tau^{-1}(\tau^{-1/2}\|\Sigma_2^{1/2}(\hat{\beta} - \beta_0)\| - \sigma_x \sigma_\beta) \leq \frac{C_\alpha}{2\sigma_\beta} - \epsilon. \tag{C24}
\]

Define \( w = \tau^{-3/2}\Sigma_2^{1/2}(\beta - \beta_0) \). Using this, we can rewrite (3) as the following problem:

\[
\hat{w} = \arg \min_{\hat{w}} \frac{c_n}{n} \|\tau^{1/2}\Sigma_1^{1/2} Tw - \tau^{-1}\varepsilon\|^2 + \frac{c_n \tau^{-1/2} \lambda_n}{\sqrt{n}} \|\Sigma_2^{-1/2}w + \tau^{-3/2} \beta_0\|_1 - \frac{c_n \tau^{-2}}{n} \|\varepsilon\|^2 - C^\phi_n.
\]

Define \( S^w_0 = \{ w \mid c_n \tau^{-1} \sigma_x \sigma_\beta + c_\alpha/4\sigma_\beta \leq c_n \|w\| \leq c_n \tau^{-1} \sigma_x \sigma_\beta + C_\alpha/\sigma_\beta \} \). Analogous to the result proved by Lemma 13, if the solution \( \hat{w}^B \) to the following problem
the proof. Let \( \hat{w} \) satisfies Eq. (C26). Next, we construct an auxiliary optimization problem:

\[
\min_{w \in S_w^n} \frac{c_n}{n} \| \tau^{1/2} \Sigma_1^{1/2} Z \hat{w} - \tau^{1/2} \| \| 2 + \frac{c_n \tau^{-1/2} \lambda}{\sqrt{n}} \| \Sigma_2^{-1/2} w + \tau^{-1/2} \beta_0 \| 1 - \frac{c_n \tau^{-2}}{n} \| \varepsilon \| 2 - C_n^\phi \tag{C25}
\]

satisfies \( c_n \| \hat{w} \| - c_n \tau^{-1} \sigma_x \sigma_\beta \in [c_n/2\sigma_\beta + \epsilon, C_n/2\sigma_\beta - \epsilon] \) w.a.p.1, then the same holds true for \( \hat{w} \), which leads to the desired result, (C25). In light of this, without ambiguity we now directly focus on (C25), and refer to \( \hat{w} \) as \( \hat{w} \) for ease of notation.

Note that for any vector \( x \), it holds that \( \| x \| 2 = \max_u \sqrt{n} u^\top x - n \| u \| 2/4 \), and \( \| x \| 1 = \max_{\| v \| \leq 1} v^\top x \). By applying these equations to \( \| \tau^{1/2} \Sigma_1^{1/2} Z \hat{w} - \tau^{1/2} \| \| 2 \) and \( \| \Sigma_2^{-1/2} w + \tau^{-1/2} \beta_0 \| 1 \), and letting \( \tilde{u} := \Sigma_1^{1/2} u \), the problem (C25) can be reformulated as:

\[
\min_{w \in S_w^n} \max_{u \in S_u^n} \frac{c_n \tau^{1/2}}{\sqrt{n}} \tilde{u}^\top Z w - \frac{c_n \tau^{-1}}{\sqrt{n}} \tilde{u}^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n \| \Sigma_1^{-1/2} \tilde{u} \| 2}{4} + \frac{c_n \tau^{-2} \lambda_n}{\sqrt{n}} v^\top \beta_0
\]

\[
+ \frac{c_n \tau^{-1/2} \lambda_n}{\sqrt{n}} v^\top \Sigma_2^{-1/2} w - \frac{c_n \tau^{-2}}{n} \| \varepsilon \| 2 - C_n^\phi. \tag{C26}
\]

For convenience, we shall continue to employ \( u \) in place of \( \tilde{u} \) throughout the remainder of the proof. Let \( S_u^n = \{ u \| u \| \leq 4 \tau^{-1} \sqrt{C_1 C_\varepsilon} \} \). Similar to the proof of Theorem 2, w.a.p.1, the optimization problem below

\[
\min_{w \in S_w^n} \max_{u \in S_u^n} \frac{c_n \tau^{1/2}}{\sqrt{n}} u^\top Z w - \frac{c_n \tau^{-1}}{\sqrt{n}} u^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n \| \Sigma_1^{-1/2} u \| 2}{4} + \frac{c_n \tau^{-2} \lambda_n}{\sqrt{n}} v^\top \beta_0
\]

\[
+ \frac{c_n \tau^{-1/2} \lambda_n}{\sqrt{n}} v^\top \Sigma_2^{-1/2} w - \frac{c_n \tau^{-2}}{n} \| \varepsilon \| 2 - C_n^\phi \tag{C27}
\]

is equivalent to Eq. (C26). Next, we construct an auxiliary optimization problem:

\[
\phi(g, h) = \max_{0 \leq \delta \leq 4 \tau^{-1} \sqrt{C_1 C_\varepsilon}} \min_{w \in S_w^n} \frac{c_n \tau^{1/2}}{\sqrt{n}} \| g^\top u - h \| 2 + \frac{c_n \tau^{-1}}{\sqrt{n}} \| \hat{w}^\top w - \frac{c_n \tau^{-1}}{\sqrt{n}} \Sigma_1^{-1/2} \| 2 - \frac{c_n \| \Sigma_1^{-1/2} \| 2}{4}
\]

\[
+ \frac{c_n \tau^{-2} \lambda_n}{\sqrt{n}} v^\top \beta_0 + \frac{c_n \tau^{-1/2} \lambda_n}{\sqrt{n}} v^\top \Sigma_2^{-1/2} w - \frac{c_n \tau^{-2}}{n} \| \varepsilon \| 2 - C_n^\phi, \tag{C28}
\]

and both \( g \in \mathbb{R}^n \) and \( h \in \mathbb{R}^p \) are standard Gaussian vectors, independent of all other random variables. Moreover, let \( \tilde{S}_n := \{ w \| c_n/2 \sigma_\beta + \epsilon < c_n \| w \| - c_n \tau^{-1} \sigma_x \sigma_\beta < C_n/2 \sigma_\beta - \epsilon \} \), define \( \phi_{\tilde{S}_n}(g, h) \) as the optimal value of the optimization problem (C28), with \( w \in S_w^n \cap \tilde{S}_n^c \).

Lemma 19 characterizes the limiting behavior of the optimal solution to (C27), \( \hat{w} \), and

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in turn, proves the desired (C24), under conditions pertaining to the optimization problem (C28). Therefore, we only need show that conditions outlined in Lemma 19 hold. That is, we need to prove the existence of the constants \( \tilde{\phi} < \tilde{\phi}_{\tilde{S}_h} \) such that for all \( \eta > 0 \), w.a.p.1, \( \phi(g, h) < \tilde{\phi} + \eta \) and \( \phi_{\tilde{S}_h}(g, h) > \tilde{\phi}_{\tilde{S}_h} - \eta \).

Following the same argument as in the proof of Theorem 2, after maximizing over the direction of \( u \) and minimizing over the direction of \( w \), Eq. (C28) becomes equivalent to:

\[
\max_{\alpha \in K_a} \min_{\|v\| \leq 1} \frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) (\Sigma^{-1}_1 - \mu_n(\alpha, \delta) I)^{-1} \times \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) - c_n \left\| n^{1/2} \tau^{1/2} \delta h - n^{-1/2} \tau^{-1/2} \lambda \Sigma^{-1/2}_2 v \right\| \alpha \\
+ \frac{c_n \tau^{-2} \lambda n v^\top \beta_0 - c_n \tau^{-2} n}{\sqrt{n}} \|\varepsilon\|^2 - C\phi_n,
\]

where \( K_a := \{\alpha | c_n \alpha - c_n \tau^{-1} \sigma_x \sigma_{\beta} \in [c_n/4 \sigma_{\beta}, C_{\alpha} / \sigma_{\beta}] \} \). By Lemma 15, the objective function of the above optimization problem is convex in \( \alpha \) and jointly concave in \((\delta, v)\). Consequently, we can interchange the order of min and max by applying Corollary 3.3 in Sion (1958). Applying \( \|x\| = \min_{\gamma > 0} \frac{1}{2\gamma} \|x\|^2 + \frac{\gamma}{2} \) to \( \left\| n^{-1/2} \tau^{-1/2} \delta h - n^{-1/2} \tau^{-1/2} \lambda \Sigma^{-1/2}_2 \right\| \alpha \), we obtain:

\[
\min_{\alpha \in K_a} \max_{\|v\| \leq 1} \max_{\delta \in [c_n/4 \tau^{-1} \sqrt{\Sigma} \gamma, 1/2]} \min_{\gamma > 0} \left\{ -\frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) (\Sigma^{-1}_1 - \mu_n(\alpha, \delta) I)^{-1} \times \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) - \frac{c_n \gamma}{2} + \frac{c_n \tau^{-2} \lambda n}{\sqrt{n}} v^\top \beta_0 - \frac{c_n \tau^{-2}}{n} \|\varepsilon\|^2 \right. \\
- \left. \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \delta h - n^{-1/2} \tau^{-1/2} \lambda \Sigma^{-1/2}_2 \right\|^2 - C\phi \right\}.
\]

By completing the square for terms associated with \( v \), we can rewrite this problem as:

\[
\min_{\alpha \in K_a} \max_{\|v\| \leq 1} \min_{\gamma > 0} \left\{ -\frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) (\Sigma^{-1}_1 - \mu_n(\alpha, \delta) I)^{-1} \times \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma^{-1/2}_1 \right) - \frac{c_n \gamma}{2} + \frac{c_n \gamma \tau^{-3}}{2 \alpha^2} \beta_0 \Sigma^{-1}_2 \beta_0 + \frac{c_n \tau^{-2} \lambda n}{\sqrt{n}} h^\top \Sigma^{-1/2}_2 \beta_0 \right. \\
- \left. \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \lambda \Sigma^{-1/2}_2 \beta_0 - \frac{\gamma}{\alpha^2} \tau^{-3}/2 \Sigma^{-1/2}_2 \beta_0 \right\|^2 - \frac{c_n \tau^{-2}}{n} \|\varepsilon\|^2 - C\phi \right\}.
\]

Denote the objective function as \( Q_n(\alpha, \delta, \gamma) \). Similar to Theorem 2, we define \( \gamma = \tau^{-1} \gamma_1 \), \( \delta = \tau^{-1} \delta_1^* + \delta_2^* + c_n^{-1/2} \delta_3 \), and \( \alpha = \tau^{-1} \sigma_x \sigma_{\beta} + c_n^{-1} \alpha_2 \). We obtain the modified objective
function \( \tilde{Q}_n(\alpha_2, \delta_3, \gamma_1) \) as follows:
\[
\tilde{Q}_n(\alpha_2, \delta_3, \gamma_1) := Q_n(\tau^{-1}\sigma_x \sigma_\beta + c_n^{-1}\alpha_2, \tau^{-1}\delta_1 + \delta_2^* + c_n^{-1/2}\delta_3, \tau^{-1}\gamma_1).
\] (C30)

Note that \( \delta_3 \in K_{\delta_3} := [-c_n^{1/2}(\tau^{-1}\delta_1^* + \delta_2^*), 4c_n^{1/2}\tau^{-1}\sqrt{C_1C_\varepsilon} - c_n^{1/2}(\tau^{-1}\delta_1^* + \delta_2^*)] \). Finally, Lemma 20 verifies the following inequalities:
\[
\phi(g, h) = \min_{\alpha_2 \in [\frac{c_\alpha}{2\sigma_\beta}, \frac{c_\alpha}{\sigma_\beta}]} \max_{\delta_3 \in K_{\delta_3} > 0} \tilde{Q}_n(\alpha_2, \delta_3, \gamma_1) < -\frac{C_\lambda}{8C_2} + \eta,
\]
\[
\phi_{\tilde{S}}(g, h) = \min_{\alpha_2 \in [\frac{c_\alpha}{2\sigma_\beta}, \frac{c_\alpha}{\sigma_\beta} + \epsilon] \cup [\frac{c_\alpha}{2\sigma_\beta} - \epsilon, \frac{c_\alpha}{\sigma_\beta}]} \max_{\delta_3 \in K_{\delta_3}} \tilde{Q}_n(\alpha_2, \delta_3, \gamma_1) > -\frac{C_\lambda}{100C_2} - \eta,
\] (C31)
which hold for sufficiently small \( \epsilon > 0 \) and \( \eta > 0 \). With Lemma 19, we conclude the proof.

**C.5 Proof of Proposition 1**

*Proof.* Since the out-of-sample data are mutually independent, Lemmas 2 and 3 lead to:
\[
\sum_{i \in OOS} y_i^2 = n_{oos}(\sigma_\varepsilon^2 + \tau \sigma_x \sigma_\beta^2) + O_P(n_{oos}^{1/2}).
\]

By the same argument, we obtain
\[
\sum_{i \in OOS} y_i^2 - (y_i - X_i \hat{\beta}_r(\lambda_n^{opt}))^2 = n_{oos}p^{-1}n\tau^2 \Delta(\hat{\beta}_r(\lambda_n^{opt})) + O_P(n_{oos}^{1/2}\tau).
\]

Theorem 2 has established \( \Delta(\hat{\beta}_r(\lambda_n^{opt})) = -\theta \sigma_x^4 \sigma_\beta^4 \sigma_\varepsilon^{-2} + o_P(1) \). Thus, as \( n_{oos}p^{-2}n^2\tau^2 \rightarrow 0 \),
\[
\sum_{i \in OOS} y_i^2 - (y_i - X_i \hat{\beta}_r(\lambda_n^{opt}))^2 = n_{oos}p^{-1}n\tau^2 \theta \sigma_x^4 \sigma_\beta^4 \sigma_\varepsilon^{-2}(1 + o_P(1)).
\]

The estimates above provide the key components for deriving the limit of \( R_{oos}^2 \):
\[
R_{oos}^2 = \frac{\sum_{i \in OOS} y_i^2 - (y_i - X_i \hat{\beta}_r(\lambda_n^{opt}))^2}{\sum_{i \in OOS} y_i^2} = p^{-1}n\theta_2(R^2)^{(1 + o_P(1))},
\]
which concludes the proof. \( \square \)
C.6 Proof of Theorem 5

Proof. For convenience, let $\hat{\beta} := \hat{\beta}_r(\lambda_n)$. We write the prediction error of the benchmark as:

$$y_{\text{new}} - \hat{y}_{b_{\text{new}}} = (w_{\text{new}})^T \gamma_0 + (x_{\text{new}})^T \beta_0 + \varepsilon_{\text{new}} - (w_{\text{new}})^T (W^T W)^{-1} W^T (W \gamma_0 + X \beta_0 + \varepsilon)$$

$$= ((u_{\text{new}})^T - (w_{\text{new}})^T (W^T W)^{-1} W^T U) \beta_0 + (\varepsilon_{\text{new}} - (w_{\text{new}})^T (W^T W)^{-1} W^T \varepsilon).$$

Similarly, for the Ridge estimator, we have

$$y_{\text{new}} - \hat{y}_{\text{new}} = ((u_{\text{new}})^T - (w_{\text{new}})^T (W^T W)^{-1} W^T U) (\beta_0 - \hat{\beta}) + (\varepsilon_{\text{new}} - (w_{\text{new}})^T (W^T W)^{-1} W^T \varepsilon).$$

As a result, with simple algebra we obtain

$$\begin{align*}
\mathbb{E} [(y_{\text{new}} - \hat{y}_{\text{new}})^2 | \mathcal{I}] &= \mathbb{E} [(y_{\text{new}} - \hat{y}_{\text{b}_{\text{new}}})^2 | \mathcal{I}] \\
&= \mathbb{E} \left[ \left( (u_{\text{new}})^T - (w_{\text{new}})^T (W^T W)^{-1} W^T U) (\beta_0 - \hat{\beta}) \right)^2 | \mathcal{I} \right] \\
&\quad - \mathbb{E} \left[ (u_{\text{new}})^T - (w_{\text{new}})^T (W^T W)^{-1} W^T U) \beta_0 \right]^2 | \mathcal{I} \\
&\quad - 2 \mathbb{E} \left[ (u_{\text{new}})^T - (w_{\text{new}})^T (W^T W)^{-1} W^T U) \hat{\beta} (\varepsilon_{\text{new}} - (w_{\text{new}})^T (W^T W)^{-1} W^T \varepsilon) \right] | \mathcal{I} \\
&:= S_1 + S_2 + S_3.
\end{align*}$$

Below we analyze $S_1$ to $S_3$ one by one. We start with $S_2$. Using the independence of $w_{\text{new}}$ with $W, U, \mathcal{I},$ and $\beta_0$, the fact that $w_{\text{new}}$ has bounded variance, we have

$$\begin{align*}
\mathbb{E} [((w_{\text{new}})^T (W^T W)^{-1} W^T U \beta_0)^2 | \mathcal{I}] &\asymp \|(W^T W)^{-1} W^T U \beta_0\|^2 = \|(W^T W)^{-1} W^T (\Sigma_1^{1/2} \Sigma_2^{1/2} \beta_0)\|^2.
\end{align*}$$

Moreover, let $x \in \mathbb{R}^n$ be a standard Gaussian vector, independent of $(W, Z, \beta_0)$. Since $Z \Sigma_2^{1/2} \beta_0$ and $x \Sigma_2^{1/2} \beta_0$ share the same distribution, $\|(W^T W)^{-1} W^T \Sigma_1^{1/2} x\|^2 \asymp_p \|(W^T W)^{-1} W^T \Sigma_1^{1/2}\|^2_F$ by Lemma 2, $\text{Tr}(W^T W)^{-1} = o_P(p^{-1} n\tau)$, and that $\|\Sigma_2^{1/2} \beta_0\|^2 \asymp_p \tau$, it follows that

$$\|(W^T W)^{-1} W^T \Sigma_1^{1/2} Z \Sigma_2^{1/2} \beta_0\|^2 \asymp_p \|\Sigma_2^{1/2} \beta_0\|^2 \|(W^T W)^{-1} W^T \Sigma_1^{1/2}\|^2_F \asymp_p o_P(p^{-1} n\tau^2). \quad (C32)$$

Additionally, given that $u_{\text{new}}$ is mean zero, independent of the remaining terms, we have

$$\begin{align*}
\mathbb{E} [(w_{\text{new}})^T (W^T W)^{-1} W^T U \beta_0 (u_{\text{new}})^T \beta_0 | \mathcal{I}] &= 0.
\end{align*}$$

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Therefore, we have established that
\[
S_2 = \mathbb{E} \left[ ((u_{new}^\top \beta_0)^2 | \mathcal{I} \right] + o_P(p^{-1} n \tau^2) = \| \Sigma_2^{1/2} \beta_0 \|^2 + o_P(p^{-1} n \tau^2).
\]

With respect to \( S_1 \), we note that \( \hat{\beta} = n^{-1} (n^{-1} X^\top \mathcal{M}_W X + n^{-1} p \tau^{-1} \lambda \Pi)^{-1} X^\top \mathcal{M}_W (U \beta_0 + \varepsilon) \). Define \( R_X = (n^{-1} X^\top \mathcal{M}_W X + n^{-1} p \tau^{-1} \lambda \Pi)^{-1} \). By direct calculations, we have
\[
\mathbb{E} \left[ ((u_{new}^\top (W^\top W)^{-1} W^\top U \hat{\beta})^2 | \mathcal{I} \right] \leq \| (W^\top W)^{-1} W^\top U \hat{\beta} \|^2 \\
\leq 2n^{-2} \| (W^\top W)^{-1} W^\top U R_X X^\top \mathcal{M}_W U \beta_0 \|^2 + 2n^{-2} \| (W^\top W)^{-1} W^\top U R_X X^\top \mathcal{M}_W \varepsilon \|^2. \tag{C33}
\]

For the second term in (C33), we first note that for any constant \( \lambda > 0 \),
\[
\| R_X X^\top \mathcal{M}_W X R_X \| = \| R_U U^\top \mathcal{M}_W U R_U \| = \frac{n \lambda (n^{-1} U^\top \mathcal{M}_W U)}{(\lambda (n^{-1} U^\top \mathcal{M}_W U) + n^{-1} p \tau^{-1} \lambda)^2} \asymp n^{-1} p^{-1} \tau^2,
\]
since \( \| n^{-1} U^\top \mathcal{M}_W U \| \leq p^{-1} n^{-1} \| U \|^2 \). Therefore, by Lemma 6 and using inequality \( \text{Tr}(AB) \leq \| A \| \text{Tr}(B) \), for any \( A = A^\top \) and \( B \geq 0 \), we have
\[
n^{-2} \| (W^\top W)^{-1} W^\top U R_X X^\top \mathcal{M}_W \varepsilon \|^2 \\
\asymp n^{-2} \text{Tr}((W^\top W)^{-1} W^\top U R_X X^\top \mathcal{M}_W X R_X U^\top W (W^\top W)^{-1}) \\
=n^{-2} \text{Tr}(R_X X^\top \mathcal{M}_W X R_X U^\top W (W^\top W)^{-2} W^\top U) \\
\leq n^{-2} \| R_X X^\top \mathcal{M}_W X R_X \| \text{Tr}((U^\top W (W^\top W)^{-2} W^\top U) \\
\leq p^{-1} \tau^2 \text{Tr}((U^\top W (W^\top W)^{-2} W^\top U) \leq p^{-1} \tau^2 \| U^\top W \| \text{Tr}((W^\top W)^{-1}) = o_p(p^{-1} n \tau^3).
\]

Similarly, we can prove that the first term in (C33) is of order \( o_p(p^{-1} n \tau^3) \). Therefore, we have \( \mathbb{E} \left[ ((u_{new}^\top (W^\top W)^{-1} W^\top U \hat{\beta})^2 | \mathcal{I} \right] = o_P(p^{-1} n \tau^3) \). With (C32) and (C33), we have
\[
\mathbb{E} \left[ ((u_{new}^\top (W^\top W)^{-1} W^\top U (\beta_0 - \hat{\beta}))^2 | \mathcal{I} \right] \\
\leq 2 \mathbb{E} \left[ ((u_{new}^\top (W^\top W)^{-1} W^\top U (\beta_0)^2 | \mathcal{I} \right] + 2 \mathbb{E} \left[ ((u_{new}^\top (W^\top W)^{-1} W^\top U (\hat{\beta})^2 | \mathcal{I} \right] = o_p(p^{-1} n \tau^2).
\]

In addition, since \( u_{new} \) is independent of \( \mathcal{I} \) and \( u_{new} \), we have
\[
\mathbb{E} \left[ (u_{new}^\top ((\beta_0 - \hat{\beta}) (u_{new}^\top (W^\top W)^{-1} W^\top U (\beta_0 - \hat{\beta}) \mathcal{I} ] = 0.
\]

Therefore, we conclude that
Similarly, it can be shown that the second term is of order $AB$ by Lemma 2, and using the fact that $\text{Tr}(I) = 1$.

For the first term, using $In$ we have

$$S_1 = \mathbb{E} \left[ \left( (u^{new})^T (\beta_0 - \hat{\beta}) \right)^2 | \mathcal{I} \right] + o_P(p^{-1}n\tau^2) = \| \Sigma_2^{1/2} (\beta_0 - \hat{\beta}) \|^2 + o_P(p^{-1}n\tau^2).$$

Finally we bound $S_3$. Since $u^{new}$ and $\epsilon^{new}$ are mean zero, mutually independent, and independent of $\mathcal{I}$, along with Lemma 2 and Cauchy-Schwartz inequality, we have

$$S_3 = 2\mathbb{E} \left[ (w^{new})^T (W^T W)^{-1} W^T U \hat{\beta} (w^{new})^T (W^T W)^{-1} W^T \epsilon | \mathcal{I} \right]$$

$$\leq 2 \left( \mathbb{E} \left[ \left((w^{new})^T (W^T W)^{-1} W^T U \hat{\beta} \right)^2 | \mathcal{I} \right] \right)^{1/2} \left( \mathbb{E} \left[ \left((w^{new})^T (W^T W)^{-1} W^T \epsilon \right)^2 | \mathcal{I} \right] \right)^{1/2}$$

$$\leq_P o_P(p^{-1/2}n^{1/2}\tau^{3/2}) (\text{Tr}(W(W^T W)^{-2} W^T))^{1/2} = o_P(p^{-1}n\tau^2).$$

In total, we conclude that

$$S_1 + S_2 + S_3 = \| \Sigma_2^{1/2} (\beta_0 - \hat{\beta}) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 + o_P(p^{-1}n\tau^2).$$

Now we prove that,

$$pn^{-1}\tau^{-2}(\| \Sigma_2^{1/2} (\beta_0 - \hat{\beta}) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2) \stackrel{p}{\to} \alpha^*.$$  \hspace{1cm} (C34)

By Theorem 2, $\tilde{\beta} := n^{-1} R_U^T U \beta_0 + \epsilon$ satisfies (C34) with $\hat{\beta}$ being replaced by $\tilde{\beta}$, where $R_U := (n^{-1} U^T U + n^{-1} p \tau^{-1} \lambda I)^{-1}$. Given that $\| \Sigma_2^{1/2} (\beta_0 - \hat{\beta}) \|^2 = \| \Sigma_2^{1/2} (\beta_0 - \tilde{\beta}) + \Sigma_2^{1/2} (\hat{\beta} - \tilde{\beta}) \|^2$ and that $\| \Sigma_2^{1/2} (\beta_0 - \hat{\beta}) \| \approx_P \| \Sigma_2^{1/2} \beta_0 \| \approx_P \tau^{1/2}$, it is easy to verify that (C34) follows from $\| \hat{\beta} - \tilde{\beta} \|^2 = o(n^2 p^{-2}\tau^3)$. Below we show this bound is satisfied. Note that

$$\| \hat{\beta} - \tilde{\beta} \|^2 \leq \frac{2}{n^2} \| R_U U^T M_W (U \beta_0 + \epsilon) - R_U U^T (U \beta_0 + \epsilon) \|^2 + \frac{2}{n^2} \| (R_U - R_U) U^T (U \beta_0 + \epsilon) \|^2$$

$$\leq \frac{4}{n^2} \| R_U U^T M_W \epsilon - R_U U^T \epsilon \|^2 + \frac{4}{n^2} \| R_U U^T M_W U \beta_0 - R_U U^T U \beta_0 \|^2$$

$$+ \frac{4}{n^2} \| (R_U - R_U^T) U^T \epsilon \|^2 + \frac{4}{n^2} \| (R_U - R_U) U^T U \beta_0 \|^2.$$ 

For the first term, using $\| R_U \| \approx_P n p^{-1} \tau$, $\| U^T U \| \approx_P p$, and Lemma 2, we have

$$\frac{4}{n^2} \| R_U U^T M_W \epsilon - R_U U^T \epsilon \|^2 \approx_P \frac{1}{n^2} \text{Tr}((M_W U - U)R_U)$$

$$= \frac{1}{n^2} \text{Tr}((W^T W)^{-1} W^T U^T U) \leq \frac{1}{n^2} \| R_U^T \| \| U^T U \| \text{Tr}(W(W^T W)^{-1} W^T) \approx_P p^{-1} \tau^2 \text{rank}(W).$$

Similarly, it can be shown that the second term is of order $O_P(p^{-1} \tau^2 \text{rank}(W))$. In addition, by Lemma 2, and using the fact that $\text{Tr}(AB) \leq \| A \| \text{Tr}(B)$ and $\| R_U \| \approx_P n p^{-1} \tau$, we have
Lemma 2. Suppose that \( d \) is a random vector with i.i.d. entries, zero mean, and unit variance. Then, for any \( n \) such that \( \beta \) is a random vector with i.i.d. entries, zero mean, and unit variance. Then, for any \( n \) such that

\[
\frac{4}{n^2} \| (R_U - \widetilde{R}_U) U^\mathsf{T} \|_2 \geq \frac{1}{n^2} \text{Tr}(U(R_U - \widetilde{R}_U)^2 U^\mathsf{T}) \leq \frac{1}{n^2} \| U^\mathsf{T} U \| \text{Tr}((R_U - \widetilde{R}_U)^2)
\]

\[
\leq \frac{p}{n^2} \text{Tr}((R_U(\widetilde{R}_U^{-1} - R_U^{-1})) \widetilde{R}_U^2) = \frac{p}{n^2} \text{Tr}((R_U U^\mathsf{T} W(W^\mathsf{T} W)^{-1} W^\mathsf{T} U \widetilde{R}_U)^2) \leq \frac{p}{n^2} \| R_U \|_2 \| \widetilde{R}_U \|_2 \| U^\mathsf{T} U \|^2 \text{Tr}(W(W^\mathsf{T} W)^{-1} W)^2) \leq P \ p^{-1} \tau^4 \text{rank}(W).
\]

Similarly, the final term is of order \( O_P(p^{-1} \tau^4 \text{rank}(W)) \). To sum up, we have \( \| \hat{\beta} - \beta \|^2 = O(p^{-1} \tau^2 \text{rank}(W)) = o(n^2 p^{-2} \tau^2) \), since \( \text{rank}(W) = o(n^2 p^{-1} \tau) \).

\( \square \)

D Technical Lemmas and Their Proofs

For completeness, the following section introduces a collection of lemmas, including proofs for some. We start with the Convex Gaussian Min-max Theorem (CGMT), a pivotal theorem to our proof. For a detailed exposition of its proof, we direct readers to the work of Thrampoulidis et al. (2015). The CGMT pertains to the following optimization problems:

\[
\Phi(G) := \min_{w \in S_w} \max_{u \in S_u} u^\mathsf{T} G w + \psi(w, u), \quad \text{and} \quad \phi(g, h) := \min_{w \in S_w} \max_{u \in S_u} \| w \|^2 g^\mathsf{T} u - \| u \|^2 h^\mathsf{T} w + \psi(w, u),
\]

where \( G \in \mathbb{R}^{m \times n}, g \in \mathbb{R}^m, h \in \mathbb{R}^n, S_w \subset \mathbb{R}^n, S_u \subset \mathbb{R}^m, \) and \( \psi : \mathbb{R}^n \times \mathbb{R}^m \rightarrow \mathbb{R} \).

**Lemma 1 (CGMT).** Suppose that \( S_w \) and \( S_u \) are compact sets, \( \psi \) is continuous on \( S_w \times S_u \), and the entries of \( G, g, \) and \( h \) are i.i.d. Gaussian. Then we have the following inequality:

\[
P(\Phi(G) < c) \leq 2P(\phi(g, h) \leq c), \quad \forall c \in \mathbb{R}.
\]

Moreover, if \( S_w \) and \( S_u \) are convex sets, and \( \psi \) is convex-concave on \( S_w \times S_u \), then

\[
P(\Phi(G) > c) \leq 2P(\phi(g, h) \geq c), \quad \forall c \in \mathbb{R}.
\]

The next lemma follows from Lemma B.26 from Bai and Silverstein (2009), which addresses the convergence of a quadratic form concerning a random vector with i.i.d. entries.

**Lemma 2.** Suppose \( x = (x_1, \cdots, x_n)^\mathsf{T} \) is a random vector with i.i.d. entries, zero mean, unit variance, and a fourth moment. For any \( A \in \mathbb{R}^{n \times n} \) such that \( \frac{1}{n} \text{Tr}(AA^\mathsf{T}) \rightarrow 0 \), it holds that \( \frac{1}{n} x^\mathsf{T} A x \rightarrow \frac{1}{n} \text{Tr}(A) \stackrel{P}{\rightarrow} 0. \)

**Lemma 3.** Assume \( x = (x_1, \cdots, x_n)^\mathsf{T} \) and \( y = (y_1, \cdots, y_m)^\mathsf{T} \) are two independent random vectors with i.i.d. entries, zero mean, and unit variance. Then, for any \( A \in \mathbb{R}^{n \times m} \) satisfying \( \frac{1}{n} \text{Tr}(AA^\mathsf{T}) \rightarrow 0 \), it follows that \( \frac{1}{n} x^\mathsf{T} A y \stackrel{P}{\rightarrow} 0. \)
Proof. The conclusion follows from the fact that \( \frac{1}{n^2} \mathbb{E}(x^\top A y)^2 = \frac{1}{n^2} \text{Tr}(A A^\top) \to 0. \)

The following result pertains to the Neumann series. A detailed proof and further discussion are available in Meyer (2000).

**Lemma 4.** If \( A \) is a square matrix with \( \|A\| < 1 \), then \( I - A \) is nonsingular and \( (I - A)^{-1} = \sum_{k=0}^\infty A^k \). As a consequence, \( \| (I - A)^{-1} - \sum_{k=0}^\ell A^k \| \leq \sum_{k=\ell+1}^\infty \|A\|^k = \|A\|^\ell+1/(1 - \|A\|). \)

**Lemma 5.** Assume \( x = (x_1, \ldots, x_n)^\top \) and \( y = (y_1, \ldots, y_p)^\top \) are two independent random vectors with i.i.d. sub-exponential random variables with their sub-exponential norm bounded by \( K \). Then for any \( A \in \mathbb{R}^{n \times n} \) and \( B \in \mathbb{R}^{n \times p} \), there exists a constant \( c > 0 \) such that

\[
\begin{align*}
P \left( |x^\top A x - \mathbb{E}x^\top A x| \geq t \right) &\leq 2 \exp \left( -c \min \left\{ \frac{t^2}{K^4 \|A\|_F^2}, \frac{t^{1/2}}{K\|A\|^{1/2}} \right\} \right), \\
P \left( |x^\top B y| \geq t \right) &\leq 2 \exp \left( -c \min \left\{ \frac{t^2}{K^4 \|B\|_F^2}, \frac{t^{1/2}}{K\|B\|^{1/2}} \right\} \right).
\end{align*}
\]

**Proof.** Inequality (D1) is given by Proposition 1.1 presented in Götze et al. (2021) for the case of symmetric \( A \). To prove it for the asymmetric case, we use the fact that \( x^\top A x = x^\top (A + A^\top) x/2 \), so that we can apply (D1) to \( (A + A^\top)/2 \). Using triangle inequalities, we have \( \| (A + A^\top)/2 \|_F^2 \leq \|A\|_F^2 \) and \( \| (A + A^\top)/2 \|^{1/2} \leq \|A\|^{1/2} \), (D1) holds for asymmetric \( A \).

To prove (D2), let \( z = (x^\top, y^\top)^\top \) and \( C = \begin{pmatrix} 0_{n \times n} & B \\ 0_{p \times n} & 0_{p \times p} \end{pmatrix} \). Applying (D1), we obtain

\[
P \left( |x^\top B y| \geq t \right) = P \left( |z^\top C z| \geq t \right) \leq 2 \exp \left( -c \min \left\{ \frac{t^2}{K^4 \|C\|_F^2}, \frac{t^{1/2}}{K\|C\|^{1/2}} \right\} \right)
\]

\[
= 2 \exp \left( -c \min \left\{ \frac{t^2}{K^4 \|B\|_F^2}, \frac{t^{1/2}}{K\|B\|^{1/2}} \right\} \right).
\]

The next lemma is established in Bai and Silverstein (2009) and Chen and Pan (2012).

**Lemma 6.** Suppose \( Z \) is an \( n \times p \) matrix with i.i.d. Gaussian entries. Then for any positive constant \( \epsilon > 0 \), it holds that \( n^{-1} Z^\top Z \leq (1 + \epsilon)(1 + \sqrt{c_n})^2 \), w.p.a.1, for \( c_n = p/n \in [0, \infty] \).

**Lemma 7** (Convexity). Let \( O \subseteq \mathbb{R}^d \) be open and convex and \( D \) be a dense subset of \( O \). For \( \theta \in O \), both \( M_n(\theta) \) and \( M(\theta) \) are convex in \( \theta \). If \( M_n(\theta) \xrightarrow{p} M(\theta) \), for any \( \theta \in D \), then \( \sup_{\theta \in K} |M_n(\theta) - M(\theta)| \xrightarrow{p} 0 \), for any compact subset \( K \subseteq O \).

This lemma has been shown by Lemma 7.75 of Liese and Miescke (2008) and Cor. II.1 of Andersen and Gill (1982). Next, we present a min-convergence theorem for functions defined on an open set \( (0, \infty) \), as shown by Lemma 10 of Thrampoulidis et al. (2018).
Lemma 8. Consider a sequence of proper, convex stochastic functions $M_n : \mathbb{R}^+ \to \mathbb{R}$, and a deterministic function $M : \mathbb{R}^+ \to \mathbb{R}$, satisfying (a) $M_n(x) \xrightarrow{P} M(x)$, $\forall x > 0$; (b) there exists $z > 0$ such that $M(x) > \inf_{y > 0} M(y)$, $\forall x \geq z$. Then we have $\inf_{x > 0} M_n(x) \xrightarrow{P} \inf_{x > 0} M(x)$.

Relatedly, we introduce a lemma for functions on a diverging sequence of closed sets.

Lemma 9. Consider a sequence of closed intervals $\{[x_n, y_n]\}_{n=1}^{\infty}$ such that $\lim_{n \to \infty} x_n = -\infty$ and $\lim_{n \to \infty} y_n = +\infty$. Additionally, let there be a sequence of proper random and convex functions $M_n : [x_n, y_n] \to \mathbb{R}$, and a convex, continuous, and deterministic function $M : \mathbb{R} \to \mathbb{R}$ that satisfy: (a) $M_n(x) \xrightarrow{P} M(x)$ for every $x \in \mathbb{R}$; (b) there exists $z > 0$ such that $M(x) > \inf_{y \in \mathbb{R}} M(y)$ holds for all $|x| \geq z$. Then it holds that $\inf_{x \in [x_n, y_n]} M_n(x) \xrightarrow{P} \inf_{x \in \mathbb{R}} M(x)$.

Proof. For $n$ sufficiently large, $z \in [x_n, y_n]$. Assume $x^* \in [-z, z]$ minimizes $M(x)$. Assumption (b) in fact implies that $x^* \in (-z, z)$ and that $M(x^*) = \inf_{x \in \mathbb{R}} M(x)$. Consider the event $\inf_{x \in [x_n, y_n]} M_n(x) < M_n(x^*)$. Under this event, there exists $|z_n| > z$ and $z_n \in [x_n, y_n]$ such that $M_n(z_n) < M_n(x^*)$. The geometry implies that there exists $\theta_n \in (0, 1)$, such that either $z_n\theta_n + x^*(1 - \theta_n) = z$ or $z_n\theta_n + x^*(1 - \theta_n) = -z$ holds. Using convexity, we have

$$\min(M_n(z), M_n(-z)) \leq \theta_n M_n(z_n) + (1 - \theta_n) M_n(x^*) < M_n(x^*).$$

By taking limits on both sides, we have $\min(M(z), M(-z)) \leq M(x^*)$, which contradicts Assumption (b). Therefore, w.p.a.1, we have $\inf_{x \in [x_n, y_n]} M_n(x) \geq M_n(x^*)$. Furthermore, by Lemma 7, for all arbitrarily small $\epsilon > 0$, w.p.a.1, $\sup_{x \in [x_n, y_n]} |M_n(x) - M(x)| < \epsilon$. In addition, by definition, there exists a sequence of $z_n$, such that $|z_n| \leq z$ and $\inf_{|x| \leq z} M_n(x) \geq M_n(z_n) - \epsilon$. Combining these two inequalities with the fact that $M(x^*)$ minimizes $M$ on $\mathbb{R}$ leads to $\inf_{|x| \leq z} M_n(x) \geq M_n(z_n) - \epsilon \geq M(z_n) - 2\epsilon \geq M(x^*) - 2\epsilon$, w.p.a.1. On the other hand, $\inf_{|x| \leq z} M_n(x) \leq M_n(x^*) \xrightarrow{P} M(x^*)$. Since $\epsilon$ is arbitrary, we have $\inf_{|x| \leq z} M_n(x) \xrightarrow{P} M(x^*)$. Along with $\inf_{|x| > z} M_n(x) \geq M_n(x^*)$ and $M_n(x^*) \xrightarrow{P} M(x^*)$ by (a), we have

$$\inf_{x \in [x_n, y_n]} M_n(x) = \min \left(\inf_{|x| \leq z} M_n(x), \inf_{|x| > z} M_n(x)\right) \xrightarrow{P} M(x^*),$$

which completes the proof. □

Lemma 10. Suppose $X$ is a standard Gaussian random variable, then for $x > 0$,

$$\frac{\sqrt{2}}{\sqrt{\pi}} \exp \left(-\frac{x^2}{2}\right) \left(2x^{-3} - 12x^{-5} - 15x^{-7}\right) \leq \mathbb{E}(|X| - x)^2 \leq \frac{\sqrt{2}}{\sqrt{\pi}} \exp \left(-\frac{x^2}{2}\right) \left(2x^{-3} + 3x^{-5}\right).$$

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Proof. With integration by parts, we find

\[ \mathbb{E}(|X| - x)^2_+ = \sqrt{\frac{2}{\pi}} \int_x^\infty (t-x)^2 \exp\left(-\frac{t^2}{2}\right) dt = \sqrt{\frac{2}{\pi}} \left(-x \exp\left(-\frac{x^2}{2}\right) + (x^2 + 1) f_G(x) \right), \]

where \( f_G(x) = \int_x^\infty \exp(-t^2/2) dt \). Lemma 10 then follows from the tail inequality:

\[ \exp\left(-\frac{x^2}{2}\right) \left(\frac{1}{x} - \frac{1}{x^3} + \frac{3}{x^5} - \frac{15}{x^7}\right) \leq f_G(x) \leq \exp\left(-\frac{x^2}{2}\right) \left(\frac{1}{x} - \frac{1}{x^3} + \frac{3}{x^5}\right). \]

**Lemma 11.** Given that \( X \) is a standard Gaussian random variable, the following inequalities hold when \( x > 0 \) and \( x \) is sufficiently large:

\[ \mathbb{E}|X|(|X| - x)^2_+ \leq 2x \mathbb{E}(|X| - x)^2_+ \quad \text{and} \quad \mathbb{E}X^2(|X| - x)^2_+ \leq 2x^2 \mathbb{E}(|X| - x)^2_+. \]

Proof. We only present the proof for the first inequality, noting that the proof for the second inequality follows a parallel methodology. By Lemma 10, with \( x \) sufficiently large we have

\[ \mathbb{E}(|X| - x)^2_+ = \frac{2\sqrt{2}}{x^3\sqrt{\pi}} \exp\left(-\frac{x^2}{2}\right) + o \left(\frac{1}{x^3}\right) \exp\left(-\frac{x^2}{2}\right). \]

By integration by parts, we arrive at

\[ \mathbb{E}|X|(|X| - x)^2_+ = \frac{2\sqrt{2}}{x^2\sqrt{\pi}} \exp\left(-\frac{x^2}{2}\right) + o \left(\frac{1}{x^2}\right) \exp\left(-\frac{x^2}{2}\right). \]

The desired result directly follows from the above two estimates. \( \square \)

**Definition 1.** A centered random variable \( X \in SE(\nu^2, \alpha) \) with \( \nu > 0 \) and \( \alpha > 0 \), if \( \mathbb{E}e^{\lambda X} \leq e^{\lambda^2 \nu^2} \), for all \( \lambda \) such that \( |\lambda| < \alpha^{-1} \).

**Lemma 12.** Let \( \{x_k\}_{k=1}^\infty \) be a sequence of diverging positive numbers. Then as \( p \to \infty \), we have \( \mathbb{P} \left( \|\Sigma_2 b_0\|_\infty < x_p \log(p) \right) \) and \( \|\Sigma_2^{1/2} h\|_\infty < x_p \sqrt{\log(p)} \), where \( \Sigma_2 \) and \( b_0 \) are defined in Assumptions 1 and 3, respectively, and \( h \) is a \( p \)-dimensional standard Gaussian vector.

Proof. We only present the proof for the first inequality, noting that the proof for the second inequality follows similarly. Without loss of generality, assume \( b_{0,i} \in SE(1,1) \). Write the \((i,j)\)-th element of \( \Sigma_2 \) as \( \Sigma_{2,ij} \). By the properties of sub-exponential variables, we have

\[ (\Sigma_2 b_0)_i \in SE \left( \sum_{j=1}^p \Sigma_{2,ij}^2, \max_j |\Sigma_{2,ij}| \right). \]
Given that $\sum_{j=1}^p \Sigma_{2,ij}^2 = (\Sigma_{2}^2)_{ij} \leq \lambda_1(\Sigma_{2}^2) = C_2^2$ and $\max_j |\Sigma_{2,ij}| \leq C_2$, we conclude that $(\Sigma_2 b_0)_i \in \SE(C_2^2, C_2)$. The tail bound of sub-exponential variables yields

$$P (|\Sigma_2 b_0|_i > x_p \log(p)) \leq 2 \exp \left( -\frac{x_p \log(p)}{2C_2} \right).$$

Therefore, the result follows by

$$P (||\Sigma_2 b_0||_\infty > x_p \log(p)) \leq \sum_{i=1}^p P (|\Sigma_2 b_0|_i > x_p \log(p)) \leq 2p \exp \left( -\frac{x_p \log(p)}{2C_2} \right) \to 0. \quad \square$$

**Lemma 13.** Under the conditions of Theorem 2, define $S_w^n := \{w | c_n \tau^{-1} \sigma_x \sigma_\beta - K_\alpha \leq c_n ||w|| \leq c_n \tau^{-1} \sigma_x \sigma_\beta + K_\alpha \}$ for some $K_\alpha$ such that $|\alpha^*_2| < K_\alpha$. If the solution $\hat{w}^B$ to

$$\arg \min_{w \in S_w^n} c_n \frac{1}{n} \|\tau^{1/2} \Sigma_{1/2} Z w - \tau^{-1} \epsilon \|_2^2 + c_n^2 \lambda \|\Sigma_{2}^{-1/2} w + \tau^{-3/2} \beta_0 \|_2^2 - \frac{c_n \tau^{-2}}{n} \|\epsilon\|_2^2 - C_n^\phi$$

satisfies $c_n \|\hat{w}^B\| - c_n \tau^{-1} \sigma_x \sigma_\beta \to \alpha^*_2$, then the same holds true for $\hat{w}$ of Eq. (C5).

**Proof.** The proof of this lemma is almost identical to Lemma 5 of Thrampoulidis et al. (2018) and is therefore omitted here. \[ \square \]

**Lemma 14.** Let $\hat{w}$ denote an optimal solution of Eq. (C10). Regarding $\phi(g, h)$ and $\phi_{S_n^\alpha}(g, h)$, as introduced and discussed in relation to Eq. (C11), suppose there are constants $\tilde{\phi}$ and $\tilde{\phi}_{S_n^\alpha}$ with $\tilde{\phi} < \tilde{\phi}_{S_n^\alpha}$, such that for all $\eta > 0$, the following hold w.a.p.1 as $n \to \infty$: (a) $\phi(g, h) < \tilde{\phi} + \eta$, (b) $\phi_{S_n^\alpha}(g, h) > \tilde{\phi}_{S_n^\alpha} - \eta$. Under these conditions, we have $\hat{w} \in S_n$ w.p.a.1.

**Proof.** Denote $\Phi$ as the optimal value of the minimization problem in Eq. (C10), and $\Phi_{S_n^\alpha}$ as the optimal value when we impose the constraint $w \in S_w^n \cap S_n^\alpha$. It is evident that $\hat{w} \in S_n$ whenever $\Phi_{S_n^\alpha} > \Phi$. In what follows, we show the latter statement holds w.a.p.1.

We now introduce the notation:

$$\phi^P := \min_{w \in S_w^n} \max_{u \in S_u^n} \mathcal{R}_n(w, v, u), \quad \phi^D := \max_{w \in S_w^n} \min_{u \in S_u^n} \mathcal{R}_n(w, v, u),$$

where $\mathcal{R}_n(w, v, u)$ is given by Eq. (C11).

By the min-max inequality (see Lemma 36.1 in Rockafellar (1970)), it holds that

$$\phi_{S_n^\alpha}^P := \min_{w \in S_w^n \cap S_n} \max_{u \in S_u^n} \mathcal{R}_n(w, v, u) = \min_{w \in S_w^n \cap S_n} \frac{\max_{u \in S_u^n} \max_{0 \leq s \leq 4\tau^{-1} \sqrt{C_1 C_2}} \|u\|_2 \mathcal{R}_n(w, v, u)}{\alpha \leq s \leq 4\tau^{-1} \sqrt{C_1 C_2}}.$$
Proof. First, we prove the objective function is convex in $\alpha$. The objective function of Eq. (C14) is convex in $\alpha$. We revisit Eq. (C11):

$$
\max_{0 \leq \delta \leq 4r^{-1}/\sqrt{c_1c_2}} \min_{w \in S_0^w \cap S_0^h} \max_{u \in \mathcal{D}} R_n(w, v, u) = \phi_{\bar{S}_n}(g, h), \quad \text{and} \quad \phi^D = \max_{w \in S_0^w \cap S_0^h} \min_{u \in \mathcal{D}} \max_{v \in \mathcal{V}} R_n(w, v, u)
$$

Utilizing CGMT (Lemma 1), and in conjunction with Eq. (D3), we have

$$
P \left( \Phi_{\bar{S}_n} < \bar{\Phi}_{\bar{S}_n} - \frac{\kappa}{3} \right) \leq 2P \left( \phi^D < \bar{\Phi}_{\bar{S}_n} - \frac{\kappa}{3} \right) \leq 2P \left( \phi(g, h) < \bar{\Phi}_{\bar{S}_n} - \frac{\kappa}{3} \right). \quad \text{(D5)}
$$

Similarly, employing CGMT along with Eq. (D4), we deduce:

$$
P \left( \Phi > \bar{\phi} + \frac{\kappa}{3} \right) \leq 2P \left( \phi^D \geq \bar{\phi} + \frac{\kappa}{3} \right) \leq 2P \left( \phi(g, h) \geq \bar{\phi} + \frac{\kappa}{3} \right). \quad \text{(D6)}
$$

Under assumptions (a) and (b) in this lemma, the right-hand sides of Eqs. (D5) and (D6) vanish as $p \to \infty$, given the choice of $\eta = \kappa/3$ for $\kappa := \bar{\phi}_{\bar{S}_n} - \bar{\phi}$. Consequently, w.a.p.1, we have: $\Phi_{\bar{S}_n} \geq \bar{\Phi}_{\bar{S}_n} - \kappa/3 > \bar{\phi} + \kappa/3 \geq \Phi$, which concludes the proof. \hfill \square

Lemma 15. The objective function of Eq. (C14) is convex in $\alpha$ and jointly concave in $(\delta, v)$.

Proof. First, we prove the objective function is convex in $\alpha = \|w\|$. We revisit Eq. (C11):

$$
\max_{0 \leq \delta \leq 4r^{-1}/\sqrt{c_1c_2}} \min_{w \in S_0^w \cap S_0^h} \max_{u \in \mathcal{D}} c_n \tau^{1/2} \alpha g^\top u - c_n \tau^{-1} u^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n \tau^{-2}}{4} \varepsilon^2 \|w\| - c_n \tau^{-3/2} v^\top \Sigma_2^{-1/2} \beta_0 - \frac{c_n \tau^{-1}}{4} \|u\|^2.
$$

Note that the term $f(\alpha, u) := c_n \tau^{1/2} \alpha g^\top u - c_n \tau^{-1} u^\top \Sigma_1^{-1/2} \varepsilon - \frac{c_n \tau^{-2}}{4} \varepsilon^2 \|w\|$ is convex in $\alpha$. After maximizing over the direction of $u$, the term remains convex in $\alpha$ since

$$
\max_{\|u\|=\delta} f(\theta \alpha_1 + (1-\theta) \alpha_2, u) \leq \max_{\|u\|=\delta} \{ \theta f(\alpha_1, u) + (1-\theta) f(\alpha_2, u) \} \leq \theta \max_{\|u\|=\delta} f(\alpha_1, u) + (1-\theta) \max_{\|u\|=\delta} f(\alpha_2, u), \quad \text{for } \theta \in (0, 1).
$$

Note that from Eq. (C12), max_{\|\xi\|=\delta} f(\alpha, u) = -\frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + c_n \tau^{-3/2} v^\top \Sigma_2^{-1/2} \beta_0 - \frac{c_n \tau^{-1}}{4} \|\xi\|^2.

Next, we demonstrate that this function is jointly concave in $(\delta, v)$. It is easy to verify that $-\|c_n \lambda v - n^{-1/2} \tau^{1/2} \delta h\| \lambda$ is jointly concave in $(\delta, v)$, since $\alpha \geq 0$. Moreover, $\lambda \tau^{-3/2} v^\top \Sigma_2^{-1/2} \beta_0 - \lambda \|\Sigma_2^{-1/2} v\|^2/4$ is concave in $v$. Therefore, it suffices to prove
Lemma 16. Let $\tilde{Q}_n = \tilde{Q}(\alpha, \delta, \gamma_x)$ be defined in Eq. (C16). It holds that

\[
\tilde{Q}_n \xrightarrow{p} \tilde{Q}(\alpha, \delta, \gamma_x) := -\frac{\delta^2}{4} \theta_1 + 2\sigma_x \sigma_\beta \alpha_2 - \frac{\gamma_x^2}{4 \sigma_x^2 \sigma_\beta^2} \lambda \theta_2 - \frac{\gamma_1 \alpha_2}{\sigma_x \sigma_\beta} + \frac{(\delta^2 \gamma_1)}{8 \lambda^2 \sigma_x^2 \sigma_\beta^2} \theta_2.
\]

In addition, for any $\eta > 0$, w.p.a.1, we have

\[
\tilde{Q}_n \xrightarrow{p} \tilde{Q}(\alpha, \delta, \gamma_x).
\]
(i) $\phi(g, h) < \min_{\alpha_2 \in [-K_\alpha, K_\alpha]} \max_{\gamma_2 \geq 0} \mathcal{Q}(\alpha_2, \delta_3, \gamma_1) + \eta$;

(ii) $\phi_{\tilde{g}_n}(g, h) > \min_{\alpha_2 \in [-K_\alpha, \alpha_2^* - \epsilon]} \max_{\gamma_2 > 0} \mathcal{Q}(\alpha_2, \delta_3, \gamma_1) - \eta$;

(iii) $\min_{\alpha_2 \in [-K_\alpha, K_\alpha]} \max_{\delta_2, \gamma_2 \geq 0} \mathcal{Q}(\alpha_2, \delta_3, \gamma_1) < \min_{\alpha_2 \in [-K_\alpha, \alpha_2^* - \epsilon]} \max_{\delta_2, \gamma_2} \mathcal{Q}(\alpha_2, \delta_3, \gamma_1)$.

Proof. The notation below is defined in the proof of Theorem 2. Let $\delta_2 = \delta_2^* + c_n^{-1/2} \delta_3$. First, we demonstrate that $c_n \tau^{-1} \mu_n(\alpha, \delta) - c_n \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \xrightarrow{P} 0$. Let $f(x) := \frac{1}{n}(\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon)^\top (\Sigma_1^{-1} - x) - 2(\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon)$. Recall that $\mu_n(\alpha, \delta)$ is the solution to $f(x) = \delta^2/4$. Note that $f(x)$ exhibits a monotonic increase in $x$ when $x \leq 1/C_1$. Therefore, it suffices to show that, given any arbitrarily small $\epsilon > 0$, w.p.a.1, the following inequalities hold: $c_n \tau f(\tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + \tau c_n^{-1} \epsilon) - c_n \delta^2 \tau/4 > c_+ > 0$ and $c_n \tau f(\tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) - \tau c_n^{-1} \epsilon) - c_n \delta^2 \tau/4 < c_- < 0$, for some constants $c_+$ and $c_-$. By Lemma 2, we can deduce the following equations:

$$
\frac{c_n \tau^{-1}}{n} \epsilon \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|)^{-2} \Sigma_1^{-1/2} \epsilon \\
- \frac{c_n \tau^{-1}}{n} \text{Tr} \left[ \Sigma_1^{-1/2} \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|)^{-2} \Sigma_1^{-1/2} \Sigma_1^{-1/2} \right] = O_P(c_n \tau^{-1} n^{-1/2}).
$$

$$
\frac{c_n \alpha^2}{n} \tau^{-2} \g \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|)^{-1/2} \g \\
- \frac{c_n \alpha^2}{n} \tau^{-2} \text{Tr} \left[ \Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|^2 \right] = O_P(c_n n^{-1/2}).
$$

Additionally, by Lemma 3, we have

$$
\frac{c_n \tau^{-1}}{n} \epsilon \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|)^{-2} \g = O_P(c_n \tau^{-1/2} n^{-1/2}).
$$

Therefore, using the definition of $f(\cdot)$ we can deduce that:

$$
c_n \tau f(\tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + \tau c_n^{-1} \epsilon) \\
- \frac{c_n \tau^{-1}}{n} \text{Tr} \left[ \Sigma_1^{-1/2} \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|)^{-2} \Sigma_1^{-1/2} \Sigma_1^{-1/2} \right] \\
- \frac{c_n \alpha^2}{n} \tau^{-2} \text{Tr} \left[ \Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \| - \tau c_n^{-1} \epsilon \|^2 \right] = O_P(c_n \tau^{-1} n^{-1/2}) = o_P(1). \quad (D11)
$$

Note that for sufficiently small $x$ such that $x \| \Sigma_1 \| < 1$, 

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\[
\frac{\tau^{-1}}{n} \text{Tr} \left[ \Sigma_\varepsilon^{1/2} \Sigma_1^{-1/2} (\Sigma_1^{-1} - x\mathbb{I})^{-2} \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} - \Sigma_\varepsilon^{1/2} \Sigma_1^{1/2} (\mathbb{I} + 2x\Sigma_1) \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} \right]
= \frac{\tau^{-1}}{n} \text{Tr} \left[ \Sigma_\varepsilon^{1/2} \Sigma_1^{-1/2} (\mathbb{I} - x\Sigma_1)^{-2} \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} - \Sigma_\varepsilon^{1/2} \Sigma_1^{1/2} (\mathbb{I} + 2x\Sigma_1) \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} \right]
\leq \tau^{-1} C_1 \| (\mathbb{I} - x\Sigma_1)^{-2} - (\mathbb{I} + 2x\Sigma_1) \| \lesssim \tau^{-1} x^2,
\]
where we apply Lemma 4 in the last inequality. As a consequence, we have:

\[
\frac{\tau^{-1}}{n} \text{Tr} \left[ \Sigma_\varepsilon^{1/2} \Sigma_1^{-1/2} (\Sigma_1^{-1} - \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2)\mathbb{I} - \tau c_n^{-1} \epsilon \mathbb{I})^{-2} \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} \right]
= \frac{1}{n} \text{Tr} \left[ \Sigma_\varepsilon^{1/2} \Sigma_1^{-1/2} (\tau^{-1} \Sigma_1^{-2} + 2 (\mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + c_n^{-1} \epsilon) \Sigma_1^2) \Sigma_1^{-1/2} \Sigma_\varepsilon^{1/2} \right] + O(\tau)
= \tau^{-1} \sigma_\varepsilon^2 \theta_1 + 2 (\mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + c_n^{-1} \epsilon) \sigma_\varepsilon^2 \theta_3 + O(\tau) + o(c_n^{-1}),
\]
where the last equation follows by Assumption 5. By the same argument, it follows that:

\[
\frac{\alpha^2 \tau^2}{n} \text{Tr} \left[ (\Sigma_1^{-1} = \tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2)\mathbb{I} - \tau c_n^{-1} \epsilon \mathbb{I})^{-2} \right] = \sigma_x^2 \sigma_\beta^2 \theta_4 + O(\tau) + o(c_n^{-1}).
\]

Applying the above estimates to the left-hand-side of (D11), we can deduce that:

\[
c_n \tau f(\tau \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + \tau c_n^{-1} \epsilon) - \frac{c_n \delta^2 \tau}{4}
= c_n \tau^{-1} \sigma_\varepsilon^2 \theta_1 + 2c_n (\mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + c_n^{-1} \epsilon) \sigma_\varepsilon^2 \theta_3 + c_n \sigma_x^2 \sigma_\beta^2 \theta_4 - c_n \frac{\tau^{-1} (\delta_1^*)^2 + 2 \delta_1^* \delta_2}{4} + o_P(1)
= 2c_n (\mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + c_n^{-1} \epsilon) \sigma_\varepsilon^2 \theta_3 + c_n \sigma_x^2 \sigma_\beta^2 \theta_4 - \frac{c_n \delta^2 \tau}{2} + o_P(1).
\]

By the definition of \(\mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2)\), the right-hand side of the above equation is positive w.p.a.1. The proof of the other inequality is similar. Hence, we have proved

\[
c_n \tau^{-1} \mu_n(\alpha, \delta) - c_n \mu(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) \overset{P}{\longrightarrow} 0.
\]

Next, to analyze \(\bar{Q}_n\), we first investigate the limiting behavior of:

\[
-\frac{\delta^2}{4} \mu_n(\alpha, \delta) + \frac{1}{n} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)^\top \left( \Sigma_1^{-1} - \mu_n(\alpha, \delta)\mathbb{I} \right)^{-1} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right).
\]

By (D13), we have \(\|\mu_n(\alpha, \delta)\Sigma_1\| = O_P(\tau)\). Applying Lemma 4 again, we deduce:

\[
\left\| (\Sigma_1^{-1} - \mu_n(\alpha, \delta)\mathbb{I})^{-2} - \Sigma_1^{-2} - 2 \mu_n(\alpha, \delta)\Sigma_1^3 - 3 \mu_n^2(\alpha, \delta)\Sigma_1^4 \right\| \lesssim_P \tau^3
\]
\[\| (\Sigma_1 - \mu_n(\alpha, \delta))^{-1} - \Sigma_1 - \mu_n(\alpha, \delta)\Sigma_1^2 - \mu_n^2(\alpha, \delta)\Sigma_1^3 \| \preceq_{\text{P}} \tau^3.\]

Furthermore, by the fact that \[\| \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \| = O_{\text{P}}(n^{-1}) \] and Eq. (C13),

\[
\frac{\delta^2}{4} \mu_n(\alpha, \delta) = \frac{1}{n} \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)^{\top} (\Sigma_1 - \mu_n(\alpha, \delta))^{-1} \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)
\]

With a similar approach, we have

\[
\frac{1}{n} \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)^{\top} (\Sigma_1 - \mu_n(\alpha, \delta))^{-1} \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right) + O_{\text{P}}(\delta^3).
\]

As a consequence, based on Lemmas 2 and 3, as well as the definition of \(\alpha_2\) and the fact that \(c_n \tau^{-1} = \mu_n(\alpha, \delta) - c_n \mu(\alpha, \delta) \xrightarrow{\text{P}} 0\), we have:

\[
-c_n \delta^2 \mu_n(\alpha, \delta) + c_n \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)^{\top} (\Sigma_1 - \mu_n(\alpha, \delta))^{-1} \left( \tau^{1/2} \sigma g - \tau^{-1} \Sigma_1^{-1/2} \varepsilon \right)
\]

Finally, we examine the remainder term that contributes to \(\tilde{Q}_n\):

\[
\frac{c_n^2 \lambda^2}{4} \left( \tau^{-3/2} \Sigma_2^{1/2} \beta_0 + \frac{\alpha^2 \delta \tau^{1/2}}{\sqrt{n} \gamma} \right)^{\top} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right)^{-1} \left( \tau^{-3/2} \Sigma_2^{1/2} \beta_0 + \frac{\alpha^2 \delta \tau^{1/2}}{\sqrt{n} \gamma} \right) - \frac{c_n \tau^2 \delta^2}{2 \gamma n} \| h \|^2.
\]

Using Lemmas 2 and 3 and the assumptions on \(\Sigma_2\), this term converges in probability to:

\[
\frac{c_n^2 \lambda^2 \tau^{-2} \sigma_2^2}{4p} \text{Tr} \left[ \Sigma_2^{1/2} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right)^{-1} \Sigma_2^{1/2} \right] + c_n \tau \text{Tr} \left[ \frac{c_n \lambda^2 \alpha^4 \Sigma_2}{4 n^2 \gamma^2} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right)^{-1} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right) \right] - \frac{\alpha^2 \delta^2}{2 \gamma n}
\]

Finally, we examine the remainder term that contributes to \(\tilde{Q}_n\):

\[
\frac{c_n^2 \lambda^2 \tau^{-2} \sigma_2^2}{4p} \text{Tr} \left[ \Sigma_2^{1/2} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right)^{-1} \Sigma_2^{1/2} \right] + c_n \tau \text{Tr} \left[ \frac{c_n \lambda^2 \alpha^4 \Sigma_2}{4 n^2 \gamma^2} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right)^{-1} \left( \frac{\lambda}{4} \Sigma_2 + \frac{c_n \alpha^2 \lambda^2}{2 \gamma} \right) \right] - \frac{\alpha^2 \delta^2}{2 \gamma n}
\]

where we apply Lemma 4 and the same argument in proving Eq. (D12). Combining this estimate with (D14) we conclude that
\[
\hat{Q}_n = c_n \tau^{-1} \sigma_\beta^2 \sigma_\beta^2 - \tau^{-1} \left( \frac{\delta_1^2}{4\lambda} \sigma_\beta^2 c_n \right) - c_n \sigma_\beta^2 \theta_3 \mu^2(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + 2 \sigma_x \sigma_\beta \alpha_2 - \frac{\gamma_1^2}{4\sigma_\beta^2} \theta_2 - \frac{\gamma_1 \alpha_2}{\sigma_x \sigma_\beta} \\
- \frac{c_n \delta_1 \delta_2 \sigma_\beta^2}{2\lambda} + \frac{(\delta_1^2)\gamma \sigma_\beta^2}{8\lambda^2 \sigma_\beta^2} \theta_2 - C_n^\phi + o_p(1) \\
= -\frac{\delta_3^2 \theta_1}{4\theta_3} + 2 \sigma_x \sigma_\beta \alpha_2 - \frac{\gamma_1^2}{4\sigma_\beta^2} \theta_2 - \frac{\gamma_1 \alpha_2}{\sigma_x \sigma_\beta} + \frac{(\delta_1^2)\gamma \sigma_\beta^2}{8\lambda^2 \sigma_\beta^2} \theta_2 + o_p(1).
\]

We now proceed to establish Claims (i) to (ii). Let \( K_{\delta_3} \) be the interval \([-c_n^{1/2}(\tau^{-1} \delta_1^* + \delta_2^*), 4c_n^{1/2} \tau^{-1} \sqrt{C_1 C_\varepsilon} - c_n^{1/2}(\tau^{-1} \delta_1^* + \delta_2^*)] \). It is sufficient to demonstrate that, for any compact set \( A \subset [-K_\alpha, K_\alpha] \), the following equation holds:

\[
\phi_A(g, h) := \min_{\alpha_2 \in A} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3 \gamma_1) \xrightarrow{P} \min_{\alpha_2 \in A} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3, \gamma_1). \quad (D15)
\]

Based on this result, we can deduce

\[
\phi(g, h) = \phi_{[-K_\alpha, K_\alpha]}(g, h) \xrightarrow{P} \min_{\alpha_2 \in [-K_\alpha, K_\alpha]} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3, \gamma_1),
\]

\[
\phi_{S_n}(g, h) = \min\{\phi_{[-K_\alpha, \alpha_2^* - \varepsilon]}(g, h), \phi_{[\alpha_2^* + \varepsilon, K_\alpha]}(g, h)\} \xrightarrow{P} \min_{\alpha_2 \in [-K_\alpha, \alpha_2^* - \varepsilon] \cup [\alpha_2^* + \varepsilon, K_\alpha]} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3, \gamma_1),
\]

which lead to (i) and (ii).

Fix \( \alpha_2 \in A \) and \( \gamma_1 > 0 \), and observe that \( \lim_{\delta_3 \to \pm \infty} \hat{Q}(\alpha_2, \delta_3, \gamma_1) \to -\infty \). By the concave version of the Lemma 9, we conclude that \( \max_{\delta_3 \in K_{\delta_3}} \hat{Q}_n(\alpha_2, \delta_3, \gamma_1) \xrightarrow{P} \max_{\delta_3 \in K_{\delta_3}} \hat{Q}(\alpha_2, \delta_3, \gamma_1) \). Since \( \hat{Q}_n \) is jointly concave in \((\delta_3, \gamma_1)\), after maximizing with respect to \( \delta_3 \), the function should remain concave in \( \gamma_1 \). Moreover, consider the following equation:

\[
\max_{\delta_3 \in \mathbb{R}} \hat{Q}(\alpha_2, \delta_3, \gamma_1) = 2 \sigma_x \sigma_\beta \alpha_2 - \frac{\gamma_1^2}{4\sigma_\beta^2} \theta_2 - \frac{\gamma_1 \alpha_2}{\sigma_x \sigma_\beta} + \frac{(\delta_1^2)\gamma \sigma_\beta^2}{8\lambda^2 \sigma_\beta^2} \theta_2.
\]

As a result, \( \lim_{\gamma_1 \to \infty} \max_{\delta_3 \in \mathbb{R}} \hat{Q}(\alpha_2, \delta_3, \gamma_1) \to -\infty \). By Lemma 8, we conclude that \( \max_{\gamma_1 > 0} \hat{Q}_n(\alpha_2, \delta_3, \gamma_1) \xrightarrow{P} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3, \gamma_1) \). Since \( \hat{Q}_n(\alpha_2, \delta_3, \gamma_1) \) is convex in \( \alpha_2 \), it should retain its convexity in \( \alpha_2 \) after being maximized with respect to \( \delta_3 \) and \( \gamma_1 \). Since the above equation holds for any \( \alpha_2 \in A \), by Lemma 7, we conclude that Eq. (D15) holds. This concludes the proof of Claims (i) and (ii).

The first-order condition implies a unique solution: \( \alpha_2^* := \arg\min_{\alpha_2} \max_{\gamma_1 > 0} \hat{Q}(\alpha_2, \delta_3, \gamma_1) \), which is given by \( \theta_2 \sigma_x^3 \left( \frac{\sigma_x \theta_1}{2 \lambda^2 \sigma_\beta} - \frac{\sigma_\beta}{x} \right) \). Thus, Claim (iii) holds true, concluding the proof. \( \square \)
Lemma 17. Under the conditions of Theorem 3, there exists a constant $\tilde{c} > 0$ that depends solely on fixed constants, such that w.p.a.1, inequality (C22) holds. In addition, as $n \to \infty$, for any given fixed $\lambda > 0$, Eq. (C23) holds.

Proof. Let us fix a constant $\tilde{c}$ such that the inequality

$$\frac{c_2^2 \sigma^2}{2K c^2} - \frac{4C_2^2 (1 + \sqrt{c_n})^2}{c_n \tilde{c}} > 100$$

remains true as $n, p \to \infty$. This is possible because $(1 + \sqrt{c_n})^2 / c_n$ is bounded as $n, p \to \infty$.

Let $S := \{\lambda_j = \epsilon + p^{-9}(j-1) : 1 \leq j \leq 1 + [p^9(\tilde{c} \tau^{-1} - \epsilon)]\}$. Given $\tau^{-1} = o(p)$, the cardinality of the set satisfies $|S| \leq p^{10}$. By definition, for any $\lambda \in [\epsilon, \tilde{c} \tau^{-1}]$, there exists a $\lambda_j \in S$ such that $|\lambda - \lambda_j| \leq p^{-9}$. By Eq. (C21), we have $|\hat{R}^K_{\text{CV}}(\lambda) - \hat{R}^K_{\text{CV}}(\lambda_j)| \leq \tilde{C} |\lambda - \lambda_j| \leq \tilde{C} p^{-9}$. Therefore, if we show that

$$\inf_{\lambda_j \in S} \left\{ \hat{R}^K_{\text{CV}}(\lambda_j) - \frac{1}{n} \|\epsilon\|^2 - \|\Sigma_2^{1/2} \beta_0\|^2 \right\} > np^{-1} \tau^2$$

holds w.p.a.1, we have

$$\inf_{\lambda \in [\epsilon, \tilde{c} \tau^{-1}] - \{\lambda_{-1}\}} \left\{ \hat{R}^K_{\text{CV}}(\lambda) - \frac{1}{n} \|\epsilon\|^2 - \|\Sigma_2^{1/2} \beta_0\|^2 \right\} > np^{-1} \tau^2 - \tilde{C} p^{-9} > \frac{np^{-1} \tau^2}{2},$$

which implies Eq. (C22). By Eq. (C17), it is easy to verify that we only need to prove:

$$\inf_{\lambda_j \in S} \left\{ n^{-1} K \|Z_{(i)} \Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2 - 2 n^{-1} K \|Z_{(i)} \Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0) - \|\Sigma_2^{1/2} \beta_0\|^2 \right\} > np^{-1} \tau^2$$

holds w.p.a.1 for all $i = 1, \ldots, K$. By the independence of $Z_{(i)}$ and $\hat{\beta}_{\lambda_j}$, the first term on the left-hand-side is distributed as:

$$n^{-1} K \|Z_{(i)} \Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2 \overset{d}{=} n^{-1} K \chi^2 (K^{-1} n) \|\Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2,$$

where $\chi^2 (K^{-1} n)$ denotes a Chi-squared random variable with $K^{-1} n$ degrees of freedom. Consequently, we can deduce that:

$$P \left( \left| n^{-1} K \|Z_{(i)} \Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2 - \|\Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2 \right| \geq \frac{\log(p)}{\sqrt{n}} \|\Sigma_2^{1/2} (\hat{\beta}_{\lambda_j} - \beta_0)\|^2 \right)$$

$$= P \left( \left| n^{-1} K \chi^2 (K^{-1} n) - 1 \right| \geq \frac{\log(p)}{\sqrt{n}} \right) \leq 2 \exp(-\tilde{c}_1 \log^2(p)), \dfrac{31}{E_c}$$

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with \( \tilde{\sigma}_2 \) being another positive constant that only depends on fixed constants. For simplicity, we consolidate the constants \( \tilde{\sigma}_1 \) and \( \tilde{\sigma}_2 \) into a unified constant denoted as \( \tilde{\sigma}_1 \). By the union bound inequality, we have that with probability exceeding \( 1 - 4p^{10} \exp(-\tilde{\sigma}_1 \log^2(p)) \), the following relation holds:

\[
\inf_{\lambda_j \in \mathcal{S}} \left\{ n^{-1} K \| \zeta_{i(i)} \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - 2n^{-1} K \zeta_{i(i)} \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) - \| \Sigma_2^{1/2} \beta_0 \|^2 \right\} \\
\geq \inf_{\lambda_j \in \mathcal{S}} \left\{ \left( 1 - \frac{\log(p)}{\sqrt{n}} \right) \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \frac{\log(p)}{\sqrt{n}} \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \| - \| \Sigma_2^{1/2} \beta_0 \|^2 \right\}.
\]

Assume for now that \( \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 \geq 50np^{-1} \tau^2 \) holds. In this scenario, \( \left( 1 - \frac{\log(p)}{\sqrt{n}} \right) \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \frac{\log(p)}{\sqrt{n}} \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \| \) is monotonically increasing in \( \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \| \) since \( \frac{\log(p)}{\sqrt{n}} = o(\tau) \), hence it achieves its minimum when \( \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 = 50np^{-1} \tau^2 \). As a result, it can be shown that \( \left( 1 - \frac{\log(p)}{\sqrt{n}} \right) \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \frac{\log(p)}{\sqrt{n}} \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \| \geq n^{-1} \tau^2 \). Therefore, we only need to prove

\[
\inf_{\lambda_j \in \mathcal{S}} \left\{ \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 \right\} \geq 50np^{-1} \tau^2 \text{ holds w.p.a.1.}
\]

We now establish a uniform lower bound for \( \| \Sigma_2^{1/2} (\hat{\beta}_{j_0} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 \), which can be written as: \( \| \Sigma_2^{1/2} \hat{\beta}_{j_0} \|^2 - 2 \beta_0^T \Sigma_2 \hat{\beta}_{j_0} \). By direct calculation, we have for each \( i \),

\[
\| \Sigma_2^{1/2} \hat{\beta}_{j_0} \|^2 \geq c_2 \| \hat{\beta}_{j_0} \|^2 = c_2 \left\| \frac{1}{n} \left( X_{(-i)}^T X_{(-i)} + c_n \lambda_j \mathbb{1} \right)^{-1} X_{(-i)} y_{(-i)} \right\|^2 \\
\geq \frac{c_2}{n^2} \left\| \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j \mathbb{1} \right\|^2 \| X_{(-i)} y_{(-i)} \|^2 \geq \frac{c_2}{n^2} \left( C_2 (1 + \sqrt{c_n})^2 + c_n \lambda_j \right)^{-2} \| X_{(-i)} y_{(-i)} \|^2.
\]

Further, by Lemmas 2 and 3, we have

\[
\| X_{(-i)}^T y_{(-i)} \|^2 = \varepsilon_{(-i)}^T X_{(-i)} X_{(-i)}^T \varepsilon_{(-i)} + 2 \varepsilon_{(-i)}^T X_{(-i)} X_{(-i)}^T \beta_0 + \beta_0^T X_{(-i)}^T X_{(-i)} \beta_0 + \sigma_{\varepsilon}^2 \text{Tr}(X_{(-i)}^T X_{(-i)}) + p^{-1} \tau \sigma_{\varepsilon}^2 \text{Tr}(X_{(-i)}^T X_{(-i)} X_{(-i)}^T X_{(-i)}) + o_p(n^{-1/2}).
\]

By the fact that \( \text{Tr}(AB) \leq \| A \| \| B \| \), we have
\[ c_2 \text{Tr}(Z_{(-i)}Z_{(-i)}^\top) \leq \text{Tr}(X_{(-i)}X_{(-i)}^\top) \leq C_2 \text{Tr}(Z_{(-i)}Z_{(-i)}^\top), \]

which, along with \((np)^{-1} \text{Tr}(Z_{(-i)}Z_{(-i)}^\top) \xrightarrow{P} (K - 1)/K\) and Eq. (C18), imply that

\[ p^{-1}\tau \text{Tr}(X_{(-i)}^\top X_{(-i)}X_{(-i)}^\top X_{(-i)}) \leq p^{-1}\tau \|X_{(-i)}^\top X_{(-i)}\| \text{Tr}(X_{(-i)}^\top X_{(-i)}) \lesssim_P \tau pn = o_P(np). \]

Therefore, w.p.a.1, we obtain

\[ \frac{c_2\sigma^2_p m}{2K} \leq \|X_{(-i)}^\top y_{(-i)}\|^2 \leq 2C_2\sigma^2_p pn. \] (D20)

Consequently, uniformly over \(\lambda_j \in S\), we deduce:

\[ \|\Sigma_2^{1/2} \beta_{\lambda_j}^i\|^2 \geq \frac{c_2^2\sigma^2_p m}{2nK} (C_2(1 + \sqrt{c_n})^2 + c_n\lambda_j)^{-2}. \] (D21)

On the other hand, we have

\[ |\beta_0^\top \Sigma_2 \beta_{\lambda_j}^i| \leq \frac{1}{n} \left| \beta_0^\top \Sigma_2 \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} X_{(-i)}^\top X_{(-i)} \beta_0 \right| + \frac{1}{n} \left| \varepsilon_{(-i)}^\top X_{(-i)} \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} \Sigma_2 \beta_0 \right|. \] (D22)

To bound the first term in (D22), by Lemma 5, we deduce that, with probability exceeding \(1 - 2p^{10} \exp(-\tilde{c}_1 \log^2(p))\), we have

\[ \sup_{\lambda_j \in S} \left| \frac{1}{n} \beta_0^\top \Sigma_2 \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} X_{(-i)}^\top X_{(-i)} \beta_0 \right| - \frac{p^{-1}\tau}{n} \text{Tr} \left( \Sigma_2 \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} X_{(-i)}^\top X_{(-i)} \right) \leq p^{-1}\tau n^{1/2} \log(p). \] (D23)

Moreover, note that \(\text{Tr}(AB) \leq \|AB\| \text{rank}(AB) \leq \|A\| \|B\| \text{rank}(B)\), we have

\[ \frac{p^{-1}\tau}{n} \text{Tr} \left( \Sigma_2 \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} X_{(-i)}^\top X_{(-i)} \right) \leq \frac{p^{-1}\tau}{n} \|\Sigma_2\| \left\| \left( \frac{1}{n} X_{(-i)}^\top X_{(-i)} + c_n\lambda_j \| \right)^{-1} \frac{1}{n} X_{(-i)}^\top X_{(-i)} \right\| \text{rank}(X_{(-i)}^\top X_{(-i)}) \]

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Using the fact that \( \lambda((A + \mathbb{I})^{-1}A) = (\lambda_1(A) + 1)^{-1}\lambda_1(A) \) and that \( n^{-1}\|X_{(-i)}^TX_{(-i)}\| \leq C_2(1 + \sqrt{c_n})^2 \). Combining the above inequality with \((D23)\), we derive that, for all \( \lambda_j \in S \),

\[
\frac{1}{n} \left\| \beta_0^T \Sigma_2 \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} X_{(-i)}^T X_{(-i)} \beta_0 \right\| \leq np^{-1} \tau C_2 \frac{C_2(1 + \sqrt{c_n})^2}{C_2(1 + \sqrt{c_n})^2 + c_n \lambda_j} + p^{-1} n^{1/2} \log(p).
\]

To bound the second term in \((D22)\), we use Lemma 5. By definition, we have

\[
\frac{1}{n} \varepsilon_{(-i)}^T X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 \beta_0 = \frac{p^{-1/2} \tau^{1/2}}{n} z_{(-i)} X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 b_0.
\]

Using the fact that \( \lambda_{\min}(n^{-1}X_{(-i)}^TX_{(-i)} + c_n \lambda_j I) \geq c_n \lambda_j \geq c_n \epsilon \) and Eq. (C18), we have

\[
\left\| X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 \right\| \leq C_2 c_n^{-1} \epsilon^{-1} n^{-1/2} \|X_{(-i)}\| \lesssim np^{-1/2}.
\]

Furthermore, since \( \|A\| \leq \sqrt{\text{rank}(A)} \|A\| \), it follows that

\[
\left\| X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 \right\|_F^2 \lesssim \text{rank}(X_{(-i)}) n^2 p^{-1} \lesssim n^3 p^{-1}.
\]

Therefore, by Lemma 5, it holds that, for some constant \( \tilde{c}_1 \),

\[
P \left( \frac{1}{n} \left\| \varepsilon_{(-i)}^T X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 \beta_0 \right\| > n^{1/2} \tau^{1/2} p^{-1} \log(p) \right) \leq 2 \exp(-\tilde{c}_1 \log^2(p)).
\]

As a consequence, with probability at least \( 1 - 2p^{10} \exp(-\tilde{c}_1 \log^2(p)) \), we have

\[
\sup_{\lambda_j \in S} \frac{1}{n} \left\| \varepsilon_{(-i)}^T X_{(-i)} \left( \frac{1}{n} X_{(-i)}^T X_{(-i)} + c_n \lambda_j I \right)^{-1} \Sigma_2 \beta_0 \right\| \leq n^{1/2} \tau^{1/2} p^{-1} \log(p). \tag{D24}
\]

Therefore, taking all bounds for components of \((D22)\) altogether, we have, w.p.a.1,
\[ |\beta_0^T \Sigma_2 \hat{\beta}_i^j| \leq n p^{-1} \tau C_2^2 \frac{(1 + \sqrt{c_n})^2}{C_2 (1 + \sqrt{c_n})^2 + c_n \lambda_j} + p^{-1} \tau n^{1/2} \log(p) + n^{1/2} \tau^{1/2} p^{-1} \log(p) \]

\[ \leq 2 n p^{-1} \tau C_2^2 \frac{(1 + \sqrt{c_n})^2}{C_2 (1 + \sqrt{c_n})^2 + c_n \lambda_j}, \tag{D25} \]

for each \( \lambda_j \in S \). With (D21) and (D25), we have, w.p.a.1,

\[ \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 = \| \Sigma_2^{1/2} \beta_0 \|^2 - 2 \beta_0^T \Sigma_2 \hat{\beta}_i^j \]
\[ \geq \frac{c_2^2 \sigma_p^2}{2 n K} (C_2 (1 + \sqrt{c_n})^2 + c_n \lambda_j)^{-2} - 4 n p^{-1} \tau C_2^2 \frac{(1 + \sqrt{c_n})^2}{C_2 (1 + \sqrt{c_n})^2 + c_n \lambda_j}. \]

This inequality holds as \( n, p \to \infty \) for each \( \lambda_j \in S \). Given our initial choice for \( \tilde{c} \), it is easy to check that the right-hand side exceeds \( 50 n p^{-1} \tau^2 \), which, in turn, establishes Eq. (C22).

To prove Eq. (C23), note that

\[ \frac{1}{n} \| y(i) - X(i) \hat{\beta}^i_{r-1} \|^2 - \frac{1}{n} \| \varepsilon(i) \|^2 = \frac{1}{n} \| Z(i) \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 + \frac{2}{n} \varepsilon(i)^T Z(i) \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0). \]

By the assumption that \( Z(i) \) and \( \hat{\beta}^i_{r-1} - \beta_0 \) are independent, we have

\[ \frac{1}{n} \| Z(i) \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 = \frac{1}{n} \chi^2 (K^{-1} n) \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2. \tag{D26} \]

Using the fact that \( n^{-1} \chi^2 (K^{-1} n) = K^{-1} + O_p(n^{-1/2}) \), we have

\[ \frac{1}{n} \chi^2 (K^{-1} n) \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 = \frac{1}{K} \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 + O_p \left( \frac{1}{\sqrt{n}} \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 \right). \]

By (D26), we conclude that

\[ \frac{1}{n} \| Z(i) \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 = \frac{1}{K} \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 + O_p \left( \frac{1}{\sqrt{n}} \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 \right). \]

Additionally, by Theorem 2, we deduce:

\[ \| \Sigma_2^{1/2} (\hat{\beta}^i_{r-1} - \beta_0) \|^2 - \| \Sigma_2^{1/2} \beta_0 \|^2 = \frac{2(K - 1)}{K} n p^{-1} \tau^2 \sigma_x^4 \left( \frac{\sigma_x^2}{2 \lambda^2} - \frac{\sigma_\beta^2}{\lambda} \right) + o_p(\tau^2 n p^{-1}). \]

Hence, using the fact that \( \| \Sigma_2^{1/2} \beta_0 \| > \tau \) we derive the following equation:
\[
\frac{1}{n} \sum_{i=1}^{K} \| Z(i) \Sigma_{2}^{1/2}(\hat{\beta}_{i-1}^c - \beta_0) \|^{2} - \| \Sigma_{2}^{1/2} \beta_0 \|^{2} = \frac{2(K - 1)}{K} np^{-1} \tau^2 \theta_{2} \sigma_{x}^{4} \left( \frac{\sigma_{e}^{2}}{2\lambda^{2}} - \frac{\sigma_{\beta}^{2}}{\lambda} \right) + o_{p}(\tau^2 np^{-1}).
\]

Thus, to prove Eq. (C23), it remains to show that: \( \frac{2}{n} \varepsilon(i) Z(i) \Sigma_{2}^{1/2}(\hat{\beta}_{i-1}^c - \beta_0) = o_{p}(\tau^2 np^{-1}). \) Given the independence of \( \varepsilon(i) \) and \( Z(i) \Sigma_{2}^{1/2}(\hat{\beta}_{i-1}^c - \beta_0) \), the following result holds:

\[
\frac{2}{n} \varepsilon(i) Z(i) \Sigma_{2}^{1/2}(\hat{\beta}_{i-1}^c - \beta_0) \xrightarrow{d} \frac{2}{n} \| \Sigma_{2}^{1/2}(\hat{\beta}_{i-1}^c - \beta_0) \| \varepsilon_{(i)}^{T} x = o_{p}(\tau^2 np^{-1}),
\]

where \( x \) is a standard Gaussian vector independent of \( \varepsilon(i) \). This concludes the proof. \( \square \)

**Lemma 18.** There exists a constant \( \tilde{C}_1 \) such that, w.p.a.1, uniformly for \( \mu_1, \mu_2 \in [0, \tilde{c}^{-1}] \),

\[
pn^{-1} \tau^{-2} | R_{K-CV}(\mu_1) - R_{K-CV}(\mu_2) | \leq \tilde{C}_1 | \mu_1 - \mu_2 | + o_{p}(pn^{-1} \tau^{-2}).
\]

**Proof.** By the Woodbury identity, we deduce that

\[
\left( \frac{1}{n} X_{(-i)}^{\top} X_{(-i)} + c_n \tau^{-1} \mu^{-1} \| \right)^{-1} - c_n^{-1} \tau \mu \| = - \frac{c_n^{-2} \tau^2 \mu^2}{n} X_{(-i)}^{\top} \left( \| + \frac{c_n^{-1} \tau \mu}{n} X_{(-i)} X_{(-i)}^{\top} \right)^{-1} X_{(-i)}.
\]

Hence, we arrive at:

\[
\sup_{1 \leq i \leq K} c_n \tau^{-3} \log^{-1}(p) \left\| \left( \frac{1}{n} X_{(-i)}^{\top} X_{(-i)} + c_n \tau^{-1} \mu^{-1} \| \right)^{-1} - c_n^{-1} \tau \mu \| + \frac{c_n^{-2} \tau^2 \mu^2}{n} X_{(-i)} X_{(-i)} \right\|
\]

\[
= \sup_{1 \leq i \leq K} c_n^{-1} \mu^2 \tau^{-1} \log^{-1}(p) \left\| \frac{1}{n} X_{(-i)}^{\top} \left( \| + \frac{c_n^{-1} \tau \mu}{n} X_{(-i)} X_{(-i)}^{\top} \right)^{-1} - \| \right\| X_{(-i)}
\]

\[
\leq \sup_{1 \leq i \leq K} \mu^3 c_n^{-2} \log^{-1}(p) \left\| \frac{1}{n} X_{(-i)}^{\top} X_{(-i)} \right\|^{2} \overset{p}{\to} 0.
\]

(D27)

The last inequality is a consequence of Eq. (C18) and the fact that

\[
\left\| \left( \| + \frac{c_n^{-1} \tau \mu}{n} X_{(-i)} X_{(-i)}^{\top} \right)^{-1} - \| \right\| \leq \left\| \left( \| + \frac{c_n^{-1} \tau \mu}{n} X_{(-i)} X_{(-i)}^{\top} \right)^{-1} \right\| \cdot c_n^{-1} \tau \mu \left\| \frac{1}{n} X_{(-i)}^{\top} X_{(-i)} \right\|
\]

\[
\leq c_n^{-1} \tau \mu \left\| \frac{1}{n} X_{(-i)}^{\top} X_{(-i)} \right\|.
\]

On the other hand, by direct calculation we have:

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\begin{equation*}
\begin{aligned}
R^{K-CV}(\mu_1) - R^{K-CV}(\mu_2) \\
= & \sum_{i=1}^{K} \left( \frac{1}{n} \| X(i) \hat{\beta}_r^{i-1} \|_2^2 - \frac{1}{n} \| X(i) \hat{\beta}_r^{i-1} \|_2^2 \right) - \frac{2}{n} y(i) X(i) (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1}) \\
= & \sum_{i=1}^{K} W_{1i}(\mu_1,\mu_2) - W_{2i}(\mu_1,\mu_2).
\end{aligned}
\end{equation*}

We next investigate \( W_{1i}(\mu_1,\mu_2) \) and \( W_{2i}(\mu_1,\mu_2) \) separately. For \( W_{1i}(\mu_1,\mu_2) \), we have
\begin{equation*}
\begin{aligned}
W_{1i}(\mu_1,\mu_2) &= \frac{1}{n} (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1})^\top X(i) X(i) (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1}) + \frac{1}{n} \beta_r^{i-1} X(i) X(i) (\beta_r^{i-1} - \beta_r^{i-1}) \\
&\leq \frac{1}{n} \left\| X(i) (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1}) \right\| \cdot \left\| X(i) \hat{\beta}_r^{i-1} \right\| + \frac{1}{n} \left\| X(i) \hat{\beta}_r^{i-1} \right\| \cdot \left\| X(i) (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1}) \right\| \\
&= \left\| X(i) \hat{\beta}_r^{i-1} \right\| \cdot \left\| X(i) (\hat{\beta}_r^{i-1} - \hat{\beta}_r^{i-1}) \right\| \\
&\leq \sup_{\mu_1 \in [0,\tilde{c}^{-1}]} \frac{1}{n} \left\| X(i) \beta_r^{i-1} - X(i) \tilde{\beta}_r^{i-1} \right\| \\
&\leq \sup_{\mu_1 \in [0,\tilde{c}^{-1}]} \frac{1}{n^3/2} \left\| X(i) \right\| \left\| \left( \frac{1}{n} X(i) X(i) + c_n^{-1} \right)^{-1} - c_n^{-1} \mu_1 \right\| + \frac{c_n^{-2} \tau^2 \mu_1^2}{n} \left\| X(i) X(i) \right\| \\
&\quad \times \left\| X(i) y(i) \right\| = O_p(\tau^3 \log(p)) = o_p(c_n^{-1/2} \tau),
\end{aligned}
\end{equation*}
\begin{equation}
(D28)
\end{equation}

where we use Eq. (D27), Eq. (C18), and Eq. (D20). Additionally, it is easy to verify that
\begin{equation*}
\begin{aligned}
\sup_{\mu_1 \in [0,\tilde{c}^{-1}]} \frac{1}{n} \left\| X(i) \beta_r^{i-1} - X(i) \tilde{\beta}_r^{i-1} \right\| &= \sup_{\mu_1 \in [0,\tilde{c}^{-1}]} \frac{1}{n} \left\| X(i) \right\| \left\| \left( \frac{1}{n} X(i) X(i) + c_n^{-1} \mu_1 \right)^{-1} - c_n^{-1} \mu_1 \right\| + \frac{c_n^{-2} \tau^2 \mu_1^2}{n} \left\| X(i) X(i) \right\| \\
&\leq \frac{c_n^{-1} \tau \tilde{c}^{-1}}{n \sqrt{n}} \left\| X(i) X(i) \right\| + \frac{c_n^{-2} \tau^2 \tilde{c}^{-2}}{n^2 \sqrt{n}} \left\| X(i) X(i) X(i) y(i) \right\|.
\end{aligned}
\end{equation*}

For the first term, by Eq. (D20), we have
\begin{equation*}
\begin{aligned}
\frac{c_n^{-1} \tau \tilde{c}^{-1}}{n \sqrt{n}} \left\| X(i) X(i) \right\| &= \frac{c_n^{-1} \tau \tilde{c}^{-1}}{n \sqrt{n}} \left\| Z(i) \Sigma_2^{1/2} X(i) y(i) \right\| \\
&= \frac{d}{n \sqrt{n}} \left( \frac{c_n^{-1} \tau \tilde{c}^{-1}}{n \sqrt{n}} \right) \sqrt{\chi^2(n/K)} \left\| \Sigma_2^{1/2} X(i) y(i) \right\| \leq \frac{\tilde{C}_1}{2} c_n^{-1/2} \tau,
\end{aligned}
\end{equation*}

w.p.a.1 for some constant \( \tilde{C}_1 \) that only depends on fixed constants. The second term can be bounded in the same way. Therefore, we have
For the first term, by Lemmas 2 and 3, it is easy to verify that
\[
\sup_{\mu_1 \in [0, \tilde{c}^{-1}]} \frac{1}{\sqrt{n}} \left\| X_{(i)} \tilde{\beta}_1^{(i)} \right\|_1 \leq \tilde{C}_1 c_n^{-1/2} + o_P(c_n^{-1/2}).
\] (D29)

Analogously, we can prove that \( \frac{1}{\sqrt{n}}\left\| X_{(i)} \tilde{\beta}_1^{(i)} - \tilde{\beta}_1^{(i)} \right\|_1 \leq \tilde{C}_1 |\mu_1 - \mu_2| c_n^{-1/2} + o_P(c_n^{-1/2}) \) holds uniformly for \( \mu_1, \mu_2 \in [0, \tilde{c}^{-1}] \), where \( \tilde{C}_1 \) is a fixed constant that may vary from line to line. In light of this, we deduce that: \( \sup_{1 \leq i \leq K} W_{1i}(\mu_1, \mu_2) \leq \tilde{C}_1^2 c_n^{-1/2} |\mu_1 - \mu_2| + o_P(c_n^{-1/2}) \) holds w.p.a.1 uniformly for \( \mu_1, \mu_2 \in [0, \tilde{c}^{-1}] \).

To bound \( W_{2i}(\mu_1, \mu_2) \), we first define \( \tilde{W}_{2i}(\mu_1, \mu_2) = \frac{2}{n} y_{(i)}^T X_{(i)} \tilde{\beta}_1^{(i)} - \tilde{\beta}_1^{(i)} \). By Eq. (D28), it holds that
\[
\sup_{\mu_1, \mu_2 \in [0, \tilde{c}^{-1}]} |\tilde{W}_{2i}(\mu_1, \mu_2) - W_{2i}(\mu_1, \mu_2)| \leq \frac{2}{n} \left| y_{(i)}^T X_{(i)} \tilde{\beta}_1^{(i)} - \tilde{\beta}_1^{(i)} \right| + \frac{2}{n} \left| y_{(i)}^T X_{(i)} \tilde{\beta}_1^{(i)} - \tilde{\beta}_1^{(i)} \right| = o_P(\tau^2 \log(p)) = o_P(c_n^{-1/2}).
\]

Moreover, employing a similar argument to that used in proving Eq. (D29), we have
\[
\sup_{\mu_1, \mu_2 \in [0, \tilde{c}^{-1}]} |\tilde{W}_{2i}(\mu_1, \mu_2)| \leq |\mu_1 - \mu_2| \frac{c_n^{-1/2}}{n^2} \left| y_{(i)}^T X_{(i)} X_{(-i)} y_{(-i)} \right| + |\mu_1 - \mu_2| \frac{c_n^{-1/2}}{n^2} \left| y_{(i)}^T X_{(i)} X_{(-i)} y_{(-i)} \right|.
\]

For the first term, by Lemmas 2 and 3, it is easy to verify that
\[
\frac{c_n^{-1/2}}{n^2} \left| y_{(i)}^T X_{(i)} X_{(-i)} y_{(-i)} \right| \leq \frac{c_n^{-1/2}}{n^2} \left| \tilde{\epsilon}_{(i)}^T X_{(i)} X_{(-i)} \tilde{\epsilon}_{(-i)} \right| + \frac{c_n^{-1/2}}{n^2} \left| \tilde{\beta}_0^T X_{(i)} X_{(-i)} \tilde{\epsilon}_{(-i)} \right| + \frac{c_n^{-1/2}}{n^2} \left| \tilde{\beta}_0^T X_{(i)} X_{(-i)} \tilde{\beta}_0 \right| \leq \tilde{C}_1 c_n^{-1/2},
\]
for some constant \( \tilde{C}_1 \) w.p.a.1. The second term can be shown analogously. As a result, we have \( \sup_{1 \leq i \leq K} W_{2i}(\mu_1, \mu_2) \leq \tilde{C}_1 c_n^{-1/2} |\mu_1 - \mu_2| + o_P(c_n^{-1/2}) \) w.p.a.1, uniformly for \( \mu_1, \mu_2 \in [0, \tilde{c}^{-1}] \). Combining the bounds for \( W_{1i}(\mu_1, \mu_2) \) and \( W_{2i}(\mu_1, \mu_2) \) concludes the proof.

**Lemma 19.** Let \( \tilde{w} \) denote an optimal solution of Eq. (C27). Regarding \( \phi(g, h) \) and \( \phi_{\tilde{S}_n}(g, h) \), as introduced and discussed in relation to Eq. (C28), suppose there are constants \( \tilde{\phi} \) and \( \tilde{\phi}_{\tilde{S}_n} \) with \( \tilde{\phi} < \tilde{\phi}_{\tilde{S}_n} \), such that for all \( \eta > 0 \), the following hold w.a.p.1 as \( n \to \infty \): (a) \( \phi(g, h) < \tilde{\phi} + \eta \), (b) \( \phi_{\tilde{S}_n}(g, h) > \tilde{\phi}_{\tilde{S}_n} - \eta \). Under these conditions, we have \( \tilde{w} \in \tilde{S}_n \) w.p.a.1.
Lemma 20. There exists some sufficiently small $\epsilon > 0$, such that for any $\eta > 0$, w.p.a.1, the inequalities in (C31) hold.

Proof. By Eq. (D14) in Lemma 16, we have the following result:

$$
-\frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{1/2} \right)^\top \left( \Sigma_1^{-1} - \mu_n(\alpha, \delta) \right) \left( \tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \right)
= c_n \tau^{-1} \sigma_x^2 \beta_0^2 - c_n \sigma_x^2 \theta_3 \mu_2(\sigma_x \sigma_\beta, \delta_1^*, \delta_2) + 2 \sigma_x \sigma_\beta \alpha_2 + \frac{c_n \tau^{-2}}{n} \|\epsilon\|^2 + o_p(1).
$$

Additionally, by Lemmas 2 and 3, we deduce that

$$
\min_{\|\nu\| \leq 1} \left\{ \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \lambda_n \Sigma_2^{-1/2} v - n^{-1/2} \tau^{-1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right\|^2 \right\}.
$$

Notice that the following relationship holds:

$$
\min_{\|\nu\| \leq 1} \left\{ \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \lambda_n \Sigma_2^{-1/2} v - n^{-1/2} \tau^{-1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right\|^2 \right\}
= \frac{c_n \alpha^2}{2\gamma} \min_{\|\nu\| \leq 1} \left\{ \left( n^{-1/2} \tau^{-1/2} \lambda_n v - n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right)^\top \Sigma_2^{-1} \left( n^{-1/2} \tau^{-1/2} \lambda_n v - n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right) \right\}
\leq \frac{c_n \alpha^2}{2\gamma c_2} \min_{\|\nu\| \leq 1} \left\{ \left\| n^{-1/2} \tau^{-1/2} \lambda_n v - n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right\|^2 \right\}
= \frac{c_n \alpha^2}{2\gamma c_2} \left\| \left( n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h + \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2 \beta_0 \right) - n^{-1/2} \tau^{-1/2} \lambda_n \right\|_+^2.
$$

Similarly, the inequality can be deduced on the alternate side as well:

$$
\min_{\|\nu\| \leq 1} \left\{ \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \lambda_n \Sigma_2^{-1/2} v - n^{-1/2} \tau^{-1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right\|^2 \right\}
\geq \frac{c_n \alpha^2}{2\gamma c_2} \left\| \left( n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h + \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2 \beta_0 \right) - n^{-1/2} \tau^{-1/2} \lambda_n \right\|_+^2.
$$

Proof. The proof closely resembles that of Lemma 14 and is thus omitted.
Together with Lemma 21, we can deduce that, w.p.a.1,
\[
\min_{\|v\| \leq 1} \frac{c_n \alpha^2}{2\gamma} \left\| n^{-1/2} \tau^{-1/2} \lambda_n \Sigma_2^{-1/2} v - n^{-1/2} \tau^{1/2} \delta h - \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2^{1/2} \beta_0 \right\|^2 \in \left[ \frac{\sigma_x^2 \sigma_{\beta}^2}{4\gamma_1 C_2} C_{\lambda}, \frac{\sigma_x^2 \sigma_{\beta}^2}{\gamma_1 \epsilon_c} C_{\lambda} \right].
\]

Recall that \( \tilde{Q}_n(\alpha_2, \delta_3, \gamma_1) \) is defined in (C30). We introduce \( \tilde{Q}_n^{\text{upper}}(\alpha_2, \delta_3, \gamma_1) \), defined as:
\[
- \frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \frac{\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon}{\Sigma_1^{-1/2} \epsilon} \right)^\top \left( \Sigma_1^{-1} - \mu_n(\alpha, \delta) \right) - \frac{1}{4} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon)
\]
\[
= - \frac{c_n}{n} \gamma \sigma_x^2 \beta_0 \Sigma_2 \beta_0 + \frac{c_n}{\sqrt{n}} \gamma \tau^{-1} \delta h \Sigma_2^{1/2} \beta_0 - \frac{c_n \tau^2}{n} \epsilon^2 - C_{\phi},
\]
and \( \tilde{Q}_n^{\text{lower}}(\alpha_2, \delta_3, \gamma_1) \), defined as:
\[
- \frac{c_n \delta^2}{4} \mu_n(\alpha, \delta) + \frac{c_n}{n} \left( \frac{\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon}{\Sigma_1^{-1/2} \epsilon} \right)^\top \left( \Sigma_1^{-1} - \mu_n(\alpha, \delta) \right) - \frac{1}{4} (\tau^{1/2} \alpha g - \tau^{-1} \Sigma_1^{-1/2} \epsilon)
\]
\[
= - \frac{c_n \gamma}{2} \sigma_x^2 \sigma_{\beta}^2 C_\lambda + \frac{c_n \gamma \tau^{-3}}{\gamma_1 \epsilon c_2} \beta_0 \Sigma_2 \beta_0 + \frac{c_n \tau^{-1} \delta h}{\sqrt{n}} \Sigma_2^{1/2} \beta_0 - \frac{c_n \tau^2}{n} \epsilon^2 - C_{\phi}.
\]

Consequently, \( \tilde{Q}_n^{\text{lower}} \leq \tilde{Q}_n \leq \tilde{Q}_n^{\text{upper}} \). Note also that \( \tilde{Q}_n^{\text{lower}}(\alpha_2, \delta_3, \gamma_1) \) and \( \tilde{Q}_n^{\text{upper}}(\alpha_2, \delta_3, \gamma_1) \) maintain their convexity in \( \alpha_2 \) and joint concavity in \( (\delta_3, \gamma_1) \). By employing a similar line of reasoning as presented in Lemma 16, alongside the definitions of \( c_\alpha \) and \( C_\alpha \), it becomes evident that there exists a sufficiently small \( \epsilon > 0 \) such that
\[
\min_{\alpha_2 \in \mathbb{R}} \max_{\delta_3 \in K_{\delta_3}} \frac{\tilde{Q}_n^{\text{upper}}}{\alpha_2 \in \mathbb{R}} \max_{\delta_3 \in K_{\delta_3}} = \frac{1}{4} \delta_3 \theta_3 + 2 \sigma_x \sigma_{\beta} \alpha_2 - \frac{\gamma_{1} \alpha_2}{\sigma_x \sigma_{\beta}} C_\lambda = - \frac{C_\lambda}{8 C_2},
\]
and
\[
\min_{\alpha_2 \in \mathbb{R}} \max_{\delta_3 \in K_{\delta_3}} \frac{\tilde{Q}_n^{\text{lower}}}{\alpha_2 \in \mathbb{R}} \max_{\delta_3 \in K_{\delta_3}} = C_\lambda
\]

These results immediately yield the desired inequalities. \( \square \)

**Lemma 21.** For any \( \alpha_2, \delta_3 \in \mathbb{R} \) and \( \gamma_1 > 0 \), w.p.a.1, we have
\[
\frac{C_\lambda}{2} \leq c_n \tau^{-1} \left\| \left( \left| n^{-1/2} \tau^{-1/2} \Sigma_2^{1/2} \delta h + \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_2 \beta_0 \right| - n^{-1/2} \tau^{-1/2} \lambda_n \right) \right\|^2 \leq 2 C_\lambda. \tag{D30}
\]

**Proof.** We first establish the following:
\[
c_n n^{-1} \tau^{-2} \left\{ \left\| \left( \Sigma_2^{1/2} \delta_1^* h \right) - \lambda_n \right\|^2 - \mathbb{E} \left\| \left( \Sigma_2^{1/2} \delta_1^* h \right) - \lambda_n \right\|^2 \right\} \xrightarrow{p} 0. \tag{D31}
\]
Let \( \tilde{h} := \Sigma_2^{1/2} h \); we then have \( \tilde{h} \sim \mathcal{N}(0, \Sigma_2) \). Let us denote the \((i, j)\)-th element of \( \Sigma_2 \) as \( \Sigma_{2,ij} \), thus we have

\[
\tilde{h}_i | \tilde{h}_i \overset{d}{=} \Sigma_{2,ij} \Sigma_{2,ii}^{-1} \tilde{h}_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1,
\]

where \( g_1 \) is a standard Gaussian random variable independent of \( \tilde{h}_i \). Consequently,

\[
\begin{align*}
\text{Cov} \left( (|\delta_1^* \tilde{h}_i| - \lambda_n)^2, (|\delta_3^* \tilde{h}_j| - \lambda_n)^2 \right) \\
= \mathbb{E} \left\{ (|\delta_1^* \tilde{h}_i| - \lambda_n)^2 \mathbb{E} \left[ (|\delta_1^* \tilde{h}_j| - \lambda_n)^2 - \mathbb{E} (|\delta_1^* \tilde{h}_j| - \lambda_n)^2 | \tilde{h}_i \right] \right\} \\
= \mathbb{E} \left\{ (|\delta_1^* \tilde{h}_i| - \lambda_n)^2 \mathbb{E} \left[ (\delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} \tilde{h}_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n)^2 \\
- \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} g_2 + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n)^2 \right) + \tilde{h}_i \right\} ,
\end{align*}
\]

where \( g_2 \sim \mathcal{N}(0, \Sigma_{2,ii}) \) is independent of both \( g_1 \) and \( \tilde{h}_i \). It is straightforward to confirm that the following inequality holds true:

\[
\begin{align*}
&\mathbb{E} \left[ \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} \tilde{h}_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right)^2 \\
- \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} g_2 + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right)^2 \right] + \tilde{h}_i \\
\leq \delta_1^* \Sigma_{2,ij} \Sigma_{2,ii}^{-1} (\tilde{h}_i - g_2) \cdot \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} \tilde{h}_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right) + \\
&\quad + \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} g_2 + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right) + \tilde{h}_i .
\end{align*}
\]

Applying the Cauchy-Schwarz inequality to the above inequality yields

\[
\begin{align*}
&\mathbb{E} \left[ \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} \tilde{h}_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right)^2 \\
- \left( \delta_1^* \left( \Sigma_{2,ij} \Sigma_{2,ii}^{-1} g_2 + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2} g_1 \right) | - \lambda_n \right)^2 \right] \tilde{h}_i \\
\lesssim \left( \mathbb{E} \left( |\Sigma_{2,ij} \Sigma_{2,ii}^{-1} (\tilde{h}_i - g_2)| \tilde{h}_i \right) \right)^{1/2} \left( \mathbb{E} Y \sim \mathcal{N}(0, 1) \left( |\delta_1^* \Sigma_{2,ij} Y \right) - \lambda_n \right)^2 .
\end{align*}
\]
where the last step is due to $c_2 \leq \Sigma_{2,ii} \leq C_2$. Therefore, by Lemma 11, we have

$$\text{Cov} \left( (|\delta_1^* h_i| - \lambda_n)_+^2, (|\delta_1^* h_j| - \lambda_n)_+^2 \right) \lesssim |\Sigma_{2,ij} (\Sigma_{2,ii}^{-1} h_i + \sqrt{\Sigma_{2,jj} - \Sigma_{2,ii}^{-1} \Sigma_{2,ij}^2 g_1}) - \lambda_n)_+^2 \right) \right) \right)^{1/2}$$

$$\lesssim |\Sigma_{2,ij} (\Sigma_{2,ii}^{-1} h_i + h_i^2 + \Sigma_{2,ij}^2 h_i^2 + \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* \Sigma_{2,ij} Y| - \lambda_n)_+^2$$

$$\lesssim |\Sigma_{2,ij} (1 + \tilde{h}_i) \left( |\Sigma_{2,ij}| h_i + \sqrt{\mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* \Sigma_{2,ij} Y| - \lambda_n)_+^2 \right),$$

where $\mathbb{E}$ is the expected value. The above inequality leads to:

$$\text{Var} \left( c_n n^{-1} \tau^2 \left( \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* h_i| - \lambda_n)_+^2 \right) \right) = \sum_{i,j=1}^{p} c_n n^{-2} \tau^{-4} \text{Cov} \left( (|\delta_1^* h_i| - \lambda_n)_+^2, (|\delta_1^* h_j| - \lambda_n)_+^2 \right)$$

$$\lesssim c_n^2 n^{-2} \tau^{-4} \log(p) \sum_{i=1}^{p} \mathbb{E}(|\delta_1^* h_i| - \lambda_n)_+^2 \sum_{j=1}^{p} |\Sigma_{2,ij}| \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* \Sigma_{2,ij} Y| - \lambda_n)_+^2$$

$$+ c_n^2 n^{-2} \tau^{-4} \log(p)^2 \sum_{i=1}^{p} \mathbb{E}(|\delta_1^* h_i| - \lambda_n)_+^2 \sum_{j=1}^{p} \sum_{i,j=1}^{p} \Sigma_{2,ij}$$

$$\leq c_n^2 n^{-2} \tau^{-4} \log(p) C_2 \sum_{i=1}^{p} \mathbb{E}(|\delta_1^* h_i| - \lambda_n)_+^2 \left( \sum_{j=1}^{p} \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* \Sigma_{2,ij} Y| - \lambda_n)_+^2 \right)^{1/2}$$

$$+ c_n^2 n^{-2} \tau^{-4} \log(p)^2 C_2 \sum_{i=1}^{p} \mathbb{E}(|\delta_1^* h_i| - \lambda_n)_+^2$$

$$= O \left( c_n^{1/2} n^{-1/2} \tau^{-1} \log(p) + c_n n^{-1} \tau^{-2} \log^2(p) \right) = o_n(1),$$

where we use $\sum_{j=1}^{p} \Sigma_{2,ij}^2 \leq C_2^2$ and Cauchy–Schwartz inequality in the second step. This leads to Eq. (D31). Using the same approach, we can prove

$$\left| \mathbb{E} \left( \left( \left( \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* h| - \lambda_n + \log^{-1}(p) \right) \right)^2 \right) - \mathbb{E} \left( \left( \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* h| - \lambda_n + \log^{-1}(p) \right) \right)^2 \right| = o_p(c_n^{-1} n \tau^2),$$

$$\left| \mathbb{E} \left( \left( \left( \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* h| - \lambda_n - \log^{-1}(p) \right) \right)^2 \right) - \mathbb{E} \left( \left( \mathbb{E}_{Y \sim N(0,1)} (|\delta_1^* h| - \lambda_n - \log^{-1}(p) \right) \right)^2 \right| = o_p(c_n^{-1} n \tau^2).$$

Now we are ready to establish Eq. (D30). Note that w.p.a.1, we have
where the last inequality is given by Lemma 12 and the fact that \( \tau \log^4(p) \to 0 \). Therefore, w.p.a.1, we have

\[
\begin{align*}
&c_n \tau^{-1} \left( \left\| \left( n^{-1/2} \tau^{1/2} \Sigma_{2}^{1/2} \delta h + \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_{2} \beta_0 \right) - n^{-1/2} \tau^{-1/2} \lambda_n \right\|_+ \right)^2 \\
&\leq c_n n^{-1} \tau^{-2} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) + \tau \Sigma_{2}^{1/2} \delta_2 h \right\| + n^{1/2} \frac{\gamma}{\alpha^2} \tau^{-1} \Sigma_{2} \beta_0 - \lambda_n \right)_+ \right)^2 \\
&\leq c_n n^{-1} \tau^{-2} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n + \log^{-1}(p) \right)_+ \right)^2,
\end{align*}
\]

Similarly, it holds that w.p.a.1,

\[
\begin{align*}
&c_n \tau^{-1} \left( \left\| \left( n^{-1/2} \tau^{1/2} \Sigma_{2}^{1/2} \delta h + \frac{\gamma}{\alpha^2} \tau^{-3/2} \Sigma_{2} \beta_0 \right) - n^{-1/2} \tau^{-1/2} \lambda_n \right\|_+ \right)^2 \\
&\geq \frac{1}{2} c_n n^{-1} \tau^{-2} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n - \log^{-1}(p) \right)_+ \right)^2.
\end{align*}
\]

Finally, by Lemma 10 and the fact that \( \lambda_n = o(\log(p)) \), it is easy to verify that

\[
\frac{\mathbb{E} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n \right\|_+ \right)^2}{\mathbb{E} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n - \log^{-1}(p) \right)_+ \right)^2} \to 1 \text{ and } \frac{\mathbb{E} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n \right\|_+ \right)^2}{\mathbb{E} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n + \log^{-1}(p) \right)_+ \right)^2} \to 1.
\]

Together with the definition that \( C_\lambda = \lim_{n \to \infty} pn^{-2} \tau^{-2} \mathbb{E} \left( \left\| \left( \Sigma_{2}^{1/2} \delta_1^* h \right) - \lambda_n \right\|_+ \right)^2 \) given by Eq. (9), we conclude the proof. \( \square \)

**Lemma 22.** Let \( \Sigma_1 \) and \( \Sigma_\varepsilon \) denote the covariance matrices of stationary processes with exponentially decaying correlations. As a result, both \( \| \Sigma_1 \| \) and \( \| \Sigma_\varepsilon \| \) are bounded and the conditions in Assumption 5 pertaining to these matrices are satisfied.

**Proof.** Given the symmetric Toeplitz structure of the matrices \( \Sigma_1 \) and \( \Sigma_\varepsilon \), the elements of these matrices can be defined as \( (\Sigma_1)_{ij} = \sigma_{1,|i-j|} \) and \( (\Sigma_\varepsilon)_{ij} = \sigma_{\varepsilon,|i-j|} \), respectively. By as-
umption, the sequences \( \{\sigma_{1,k}\}_{k=1}^{\infty} \) and \( \{\sigma_{\varepsilon,k}\}_{k=1}^{\infty} \) decay exponentially, i.e., there exist positive constants \( c \) and \( C \) such that \( |\sigma_{1,k}|, |\sigma_{\varepsilon,k}| \leq C \exp(-ck) \), for all \( k \).

We first show that the eigenvalues of \( \Sigma_1 \) are bounded. The proof for \( \Sigma_{\varepsilon} \) is similar. Note that

\[
\max_{i=1,\ldots,n} \sum_{j=1}^{n} |(\Sigma_1)_{ij}| = \max_{i=1,\ldots,n} \sum_{j=1}^{n} |\sigma_{1,i-j}| \leq 2 \sum_{k=0}^{n-1} |\sigma_{1,k}|.
\]

Since \( \{\sigma_{1,k}\}_{k=1}^{\infty} \) decays exponentially, the right-hand-side is bounded, as \( n \to \infty \). The bound on \( \|\Sigma_1\| \) thus follows from Gershgorin circle theorem.

Next we establish \( \frac{1}{n} \operatorname{Tr}(\Sigma_{\varepsilon} \Sigma_1) = \sigma_{\varepsilon}^2 \theta_1 + o(c_n^{-1}\tau) \). The proofs for \( \theta_3 \) and \( \theta_4 \) follow analogously. The series \( \sigma_{1,0}\sigma_{\varepsilon,0} + 2 \sum_{i=1}^{\infty} \sigma_{1,i}\sigma_{\varepsilon,i} \) is convergent, since the series \( \{\sigma_{1,k}\}_{k=1}^{\infty} \) and \( \{\sigma_{\varepsilon,k}\}_{k=1}^{\infty} \) decay exponentially. We denote the limit as \( \sigma_{\varepsilon}^2 \theta_1 \). Moreover, given that

\[
\frac{1}{n} \left[ \operatorname{Tr}(\Sigma_{\varepsilon} \Sigma_1) - n\sigma_{\varepsilon}^2 \theta_1 \right] = \frac{1}{n} \left[ -\sum_{i=1}^{n-1} 2i\sigma_{1,i}\sigma_{\varepsilon,i} - 2n \sum_{i=n}^{\infty} \sigma_{1,i}\sigma_{\varepsilon,i} \right],
\]

and that \( \{\sigma_{1,k}\}_{k=1}^{\infty} \) and \( \{\sigma_{\varepsilon,k}\}_{k=1}^{\infty} \) decay exponentially, it follows that \( -\sum_{i=1}^{n-1} 2i\sigma_{1,i}\sigma_{\varepsilon,i} - 2n \sum_{i=n}^{\infty} \sigma_{1,i}\sigma_{\varepsilon,i} < \infty \), as \( n \to \infty \). We thereby have that

\[
\frac{1}{n} \operatorname{Tr}(\Sigma_{\varepsilon} \Sigma_1) = \sigma_{\varepsilon}^2 \theta_1 + O\left(n^{-1}\right) = \sigma_{\varepsilon}^2 \theta_1 + o(c_n^{-1}\tau).
\]

This concludes our proof.

\[\square\]

References


Chen, B. and G. Pan (2012). Convergence of the largest eigenvalue of normalized sample covariance matrices when \( p \) and \( n \) both tend to infinity with their ratio converging to zero. Bernoulli 18(4), 1405 – 1420.


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