Optimal High-Dimensional Covariance Matrix Sensing With Non-convex Sparse Learning

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Abstract—Covariance matrix plays a pivotal role in highdimensional statistics, making accurate estimation of secondorder statistics a critical task. Directly estimating the covariance matrix from samples is a common approach, yet it faces substantial challenges when data changes rapidly and the acquisition devices have limited processing power and storage capacity. A classical solution is covariance matrix sensing (CMS), which involves sensing and compressing signals and estimating the covariance matrix from measurements. To efficiently estimate the covariance matrix, a prevalent assumption is sparsity. In this paper, we explore a quadratic measurement model and propose an estimator, which adopt the least squares loss function with non-convex penalty to estimate sparse covariance matrices. It is a challenging task to analyze the statistical optimality of convergence by oracle style. Fortunately, the majorizationminimization (MM) framework offers a solution via a multistage convex relaxation approach. We clearly establish the statistical performance of all the sequential approximate solutions produced by the MM-based algorithm. The performance is supported by numerical simulation.

Index Terms—Covariance matrix sensing, quadratic measurements, majorization-minimization, positive definiteness, sparsity, non-convex statistical optimization.

I. Introduction

The covariance matrix is indispensable in statistical inference [1] and information processing [2], providing critical insights into the variability and relationship between different variables [3]. A direct method is to estimate covariance matrices from the sample data [3]. However, this approach often encounters significant challenges due to the rapidly changing data, coupled with the limited processing capability and storage capacity of sampling equipment. Thus, it is desirable to explore to recover covariance matrices from single data streams and a minimal number of storage measurements. Covariance matrix sensing (CMS, a.k.a., covariance matrix sketching) is a classical solution to address this challenging issue. It is a method employed in statistics and signal processing that involve two key processes: sensing and compressing the signals, and estimating the covariance matrix from potentially noisy measurements [4]. Specifically in this paper, we explore a quadratic measurement model as follows:

$$y_i = \boldsymbol{a}_i^{\top} \boldsymbol{\Sigma} \boldsymbol{a}_i + \eta_i, \qquad i = 1, \dots, m,$$
 (1)

where $\Sigma \in \mathbb{R}^{d \times d}$ is the covariance matrix, $\{y_i\}_{i=1}^m$ means a sequence of measurements, $a_i \in \mathbb{R}^d$ denotes the sensing

vectors, $\{\eta_i\}_{i=1}^m$ represents the noise term, m is the number of measurements.

The quadratic measurement model in the form of (1) has broad applications in many fields such as wireless communication [5], optical imaging [6], environmental monitoring [7], etc. In wireless communication, the power spectrum characterizes the distribution of power or energy of a signal across different frequencies, providing crucial insights into the signal's properties and behavior [8], [9]. Consequently, estimating the power spectrum is vital. Power spectrum estimation is a signal processing technique that enables the reconstruction of signals or images from far fewer samples than traditionally required by the Nyquist-Shannon sampling theorem, used to infer the spectral characteristics of a signal or stochastic process [10]. This technique focuses on capturing the spectral properties of random signals based on energy observations over time [5]. In this approach, a sensing vector a_i is employed to interact with the signal x_t at discrete times. The vector a_i could be implemented using techniques such as random demodulators [11]. The average energy of the signal over N observations is calculated as follows:

$$y_i = \frac{1}{N} \sum_{t=1}^{N} (\boldsymbol{a}_i^{\top} \boldsymbol{x}_t)^2 = \boldsymbol{a}_i^{\top} \boldsymbol{\Sigma}_N \boldsymbol{a}_i, \text{ for } i = 1, \dots, m,$$
 (2)

where $\Sigma_N = \frac{1}{N} \sum_{t=1}^N \boldsymbol{x}_t \boldsymbol{x}_t^{\top}$ symbolizes the sample covariance matrix. The power spectrum can then be inferred by analyzing the structure of the covariance matrix, typically through spectral analysis techniques such as Fourier transforms. This sampling model ties the observed energy directly to the power spectrum, providing a robust way to estimate the spectral distribution of the signal based purely on energy measurements, making it particularly suitable for high-frequency applications where phase measurements may be difficult to obtain [5]. In high-dimensional data processing, covariance sketching emerges as a pivotal application of the quadratic model, particularly advantageous in settings where the data dimensions and volume surpass the available computational resources [4], [12]. This technique fundamentally entails the estimation of covariance matrices through the use of compressed or "sketched" data representations. Specifically, the quadratic nature of the model is employed to transform and reduce the dimensionality

of data streams or samples, which are then used to approximate the covariance matrix efficiently. For instance, considering a sequence of high-dimensional vectors $x_t \in \mathbb{R}^d$, the process involves selecting a random sketching vector $oldsymbol{a}_{i_t}$ at each time t and computing the quadratic sketch $y_i = (\boldsymbol{a}_i^{\top} \boldsymbol{x}_t)^2$. These sketches are subsequently aggregated and normalized to form an estimate of the covariance matrix $\Sigma = \mathbb{E}(x_t x_t^{\top})$, with adjustments for error terms denoted by η_i . Such methodologies not only facilitate the processing of data with limited memory and computational capabilities but also enhance the ability to handle real-time data streams efficiently, thus exemplifying the practical applications of quadratic models in modern statistical data analysis [13]. Beyond the applications previously mentioned, the quadratic measurement model is also extensively utilized in fields such as robotics and autonomous systems [14], environmental perception [15], phaseless measurements in physics [6], graph sketching [16], and direction of arrival (DoA) estimation [17], among others. Overall, each of these applications necessitates the recovery of covariance matrices from a minimal set of quadratic measurements (1).

The CMS technique is inspired by the success of compressed sensing [18], which asserts it is possible to sense and compress signals simultaneously without lose information. To achieve it effectively, exploiting low-dimensional structures inherent in high-dimensional data is a common approach [19], with sparsity being one of the widely adopted assumption [20], i.e., most off-diagonal elements are almost zero, significantly decreasing the number of parameters that need estimation. Under approximately sparse assumption, [8] and [21] investigated the recovery of second-order statistics for nearly sparse cyclostationary signals through random linear measurements, employing ℓ_1 -minimization but without performance guarantees. Another pioneering work by Dasarathy et al. [4] have suggested estimating a nearly sparse covariance matrix via measurements formatted as $Y = B\Sigma B^{\top}$, where B represents a sketching matrix based on expander graphs.

The ℓ_1 penalty has been utilized in CMS. [22] derived a fundamental guarantee on how many samples are sufficient to approximate the ground truth by ℓ_1 regularization and showed that covariance estimation from compressive measurements can be highly robust under sparse assumption. However, it is widely acknowledge that it introduces a non-negligible bias into the estimator. To mitigate this bias, alternatives like the non-convex smoothly clipped absolute deviation (SCAD) penalty [23] and minimax concave penalty (MCP) [24] have been suggested. The benefits of non-convex penalties can also be found in many other topics such as sparse covariance estimation [25], low-rank matrix recovery [26], high-dimensional graphical models [27], graph trend filtering [28], etc.

Based on these insights, we proposed to estimate covariance matrix from a small number of measurements using the non-convex penalty. And we adopts a sub-Gaussian error assumption, different from the error term in [22] which is assumed to be bounded in ℓ_1 or ℓ_2 norm. Analyzing the statistical optimality of oracle-style convergence is a complex task due to the non-convexity of the penalty function. To address this,

we designed a multistage convex relaxation algorithm using a majorization-minimization (MM) approach that solves a series of convex subproblems. This ensures that the approximate local optimum shares the same optimal statistical properties as the unattainable global optimum. Our findings also confirm that the estimator produced by the MM-based algorithm reaches the oracle statistical rate under minimal assumptions. Theoretical analyses are supported by numerical experiments, which demonstrate the method's effectiveness.

II. PROBLEM FORMULATION

We denote the matrix to be estimated as $\Sigma \in \mathbb{R}^{d \times d}$. In general, it is NP-hard to estimate the covariance matrix $\Sigma \in \mathbb{R}^{d \times d}$ from $m < \frac{d(d+1)}{2}$ measurements, unless appropriate structural assumptions, such as sparsity, are assumed to be known as a prior. In this paper, we concentrate on the sparse covariance matrices estimation. Additionally, we outline some standard assumptions on sensing vectors and noise in Assumption 1.

Assumption 1. We operate under the assumption that $\{\eta_i\}_{i=1}^m$ and $\{a_i\}_{i=1}^m$ consist of i.i.d sub-Gaussian random variables. Specifically, η_i $(1 \le i \le m)$ is drawn from a sub-Gaussian distribution characterized by a mean of 0 and a variance proxy σ^2 . Similarly, a_i 's $(1 \le i \le m)$ are considered i.i.d copies of $z = [z_1, \dots, z_d]^\top$, where each z_i is i.i.d drawn from a distribution satisfying

$$\mathbb{E}(z_i) = 0, \mathbb{E}(z_i^2) = 1, \text{ and } \mathbb{E}(z_i^4) > 1.$$
 (3)

We propose to estimate the sparse covariance matrices from quadratic measurements using the non-convex penalty. The idea is simple yet powerful. For notational simplicity, let $\boldsymbol{A}_i \coloneqq \boldsymbol{a}_i \boldsymbol{a}_i^\top$ represent the equivalent sensing matrix, $\boldsymbol{Y} \coloneqq [y_1, \cdots, y_m]^\top$ express the equivalent measurements vector, and define the linear operator $\boldsymbol{\mathcal{A}}(\boldsymbol{\Sigma}) : \mathbb{R}^{d \times d} \mapsto \mathbb{R}^m$ that maps a matrix $\boldsymbol{\Sigma} \in \mathbb{R}^{d \times d}$ to $\{\langle \boldsymbol{\Sigma}, \boldsymbol{A}_i \rangle\}_{i=1}^m$. Thus, the measures \boldsymbol{Y} obeys $y_i \coloneqq \langle \boldsymbol{A}_i, \boldsymbol{\Sigma} \rangle + \eta_i$ and we can express the estimator as

$$\min_{\Sigma \succ \mathbf{0}} \left\{ \frac{1}{2m} \left\| \mathbf{Y} - \mathcal{A} \left(\mathbf{\Sigma} \right) \right\|_{2}^{2} - \tau \log \det \mathbf{\Sigma} + \sum_{j,k} p_{\lambda} \left(\left| \Sigma_{jk} \right| \right) \right\}. \tag{4}$$

In (4), the first term aims to minimize the least squares errors. The second term, a log-determinant barrier function, ensures positive definiteness with a barrier parameter $\tau \geq 0$. The third term p_{λ} represents a series of non-convex penalty functions governed by a regularization parameter $\lambda > 0$. We impose certain restrictions on it, as illustrated in Assumption 2.

Assumption 2. The function $p_{\lambda}(t)$ defined on $[0, +\infty)$ satisfies:

- $p_{\lambda}(t)$ is non-decreasing on $[0, +\infty)$ with $p_{\lambda}(0) = 0$ and is differentiable almost everywhere on $(0, +\infty)$;
- $0 \le p_{\lambda}'(t_1) \le p_{\lambda}'(t_1) \le \lambda$ for all $t_1 \ge t_2 \ge 0$ and $\lim_{t\to 0} p_t'(t) = \lambda$;

- There exists an $\alpha > 0$ such that $p'_{\lambda}(t) = 0$ for $t \ge \alpha \lambda$;
- There exists some $c \in (0, \alpha)$ such that $p'_{\lambda}(c\lambda) \geq \frac{\lambda}{2}$.

Under Assumption 2, the initial trio of conditions guarantee sparsity and unbiasedness. The final criterion is invariably satisfied on account of $p'_{\lambda}(0) = \lambda$ and $p'_{\lambda}(\alpha\lambda) = 0$ which is listed for the convenience of subsequent theoretical analysis. Typical examples of $p_{\lambda}(\cdot)$ in Assumption (2) include SCAD [23] and MCP [23].

III. OPTIMIZATION ALGORITHM

We now construct an MM-based algorithm to solve (4). The multistage convex relaxation is essentially a sequential optimization framework [29]. At each stage, we take weighted ℓ_1 -norm as surrogate function of $\sum_{j,k} p_{\lambda}\left(|\Sigma_{jk}|\right)$. For notational simplicity, we define $f(\Sigma) = \frac{1}{2m} \|Y - \mathcal{A}(\Sigma)\|_2^2$ $\tau \log \det \Sigma$. Consequently, we consider to solve a sequence of convex relaxation subproblems as follows:

$$\min_{\Sigma \succ \mathbf{0}} \left\{ f(\Sigma) + \sum p_{\lambda}' \left(\left| \widehat{\Sigma}_{jk}^{(q-1)} \right| \right) |\Sigma_{jk}| \right\}, \quad 1 \le q \le Q,$$
(5)

where $\widehat{oldsymbol{\Sigma}}^{(q)}$ is the optimal solution to the q-th subproblem.

Each subproblem corresponds to a weighted ℓ_1 -penalized covariance estimation problem, which can generally be reformulated as follows:

$$\min_{\mathbf{\Sigma} \subset \mathbf{0}} \left\{ f(\mathbf{\Sigma}) + \| \mathbf{\Lambda} \odot \mathbf{\Sigma} \|_{1} \right\}, \tag{6}$$

where Λ is the regularized parameter matrix with $\Lambda_{jk}=$ $p_{\lambda}'\left(\left|\widehat{\Sigma}_{jk}\right|\right) \in [0,\lambda]$. According to the Karush-Kuhn-Tucker (KKT) conditions, the unique sparse global optimal solution $\widehat{\Sigma}$ for (6) satisfies the first-order optimal condition:

$$\nabla f\left(\widehat{\boldsymbol{\Sigma}}\right) + \boldsymbol{\Lambda} \odot \widehat{\boldsymbol{\Xi}} = 0, \text{ with } \widehat{\boldsymbol{\Xi}} \in \partial \left\|\widehat{\boldsymbol{\Sigma}}\right\|_{1}$$

in which $\nabla f\left(\mathbf{\Sigma}\right)=-\frac{1}{m}\mathcal{A}^{*}\left(\mathbf{Y}-\mathcal{A}\left(\mathbf{\Sigma}\right)\right)- au\mathbf{\Sigma}^{-1}.$ Obviously, it is impossible to find an exact solution of problem (6) which has no analytical solution. In practice, we settle for the secondbest solution defined in 3 to (6), and terminate the iterations when the approximate KKT condition holds.

Definition 3. Given a tolerance level ε , $\widetilde{\Sigma}^{(q)}$ qualifies as an ε -optimal solution to (6) if $\omega_{\Lambda^{(q-1)}}\left(\widetilde{\Sigma}^{(q)}\right) \leq \varepsilon$, where

$$\omega_{\pmb{\Lambda}^{(q-1)}}\left(\widetilde{\pmb{\Sigma}}\right) = \min_{\pmb{\Xi} \in \partial \|\pmb{\Sigma}\|_1} \left\| \nabla f\left(\pmb{\Sigma}\right) + \pmb{\Lambda}^{(q-1)} \odot \pmb{\Xi} \right\|_{\max},$$

where
$$\Lambda_{jk}^{(q-1)} = p_{\lambda}' \left(\left| \widehat{\Sigma}_{jk}^{(q-1)} \right| \right)$$
.

where $\Lambda_{jk}^{(q-1)} = p_{\lambda}' \left(\left| \widehat{\Sigma}_{jk}^{(q-1)} \right| \right)$. The MM algorithm is described in Algorithm 1, where we set $\Sigma^{\{0\}}=\mathbf{I}$ as a trivial start. To obtain $\widetilde{\Sigma}^{(q)}$ by solving (6), we applying a proximal Newton method with backtracking line search [30], [31]. We start with $\Sigma_0^{(q)} = \widetilde{\Sigma}^{(q-1)}$ and stop the proximal Newton algorithm when $\omega_{\Lambda^{(q-1)}}\left(\widetilde{\Sigma}\right)$ researches the prefixed optimization error ε , which is given in Definition on this page.

Algorithm 1: The MM-Based Multistage Convex Relaxation Algorithm for Solving (4).

Input:
$$\mathcal{S}, \tau, \lambda$$
;

1 Initialize $\widetilde{\boldsymbol{\Sigma}}^{(0)} = \mathbf{I}$

2 for $q = 1, 2, \dots, Q$ do

3 $\left| \begin{array}{c} \Lambda_{ij}^{(q-1)} = p_{\lambda}' \left(\left| \widetilde{\Sigma}_{ij}^{(q-1)} \right| \right); \\ \mathbf{4} & \text{obtain } \widetilde{\boldsymbol{\Sigma}}^{(q)} \text{ by solving (6);} \\ \mathbf{5} & q = q+1; \\ \mathbf{6} \text{ end} \\ \mathbf{Output: } \widetilde{\boldsymbol{\Sigma}}^{(Q)} \end{array} \right.$

IV. STATISTICAL THEORIES

In this section, we first introduce some important definitions and technical assumptions. Following this foundational setup, we establish the statistical convergence rate of our proposed covariance estimator.

A. Assumptions

Definition 4 (Sparse Eigenvalue). Given any positive integer s, we define the largest and smallest s-sparse eigenvalues of $\nabla^2 f(\mathbf{\Sigma})$ as

$$\begin{split} \rho_{s}^{+} &= \sup \left\{ \mathbf{v}^{\top} \nabla^{2} f\left(\mathbf{\Sigma}\right) \mathbf{v} \mid \left\|\mathbf{v}\right\|_{2}^{2} = 1, \left\|\mathbf{v}\right\|_{0} \leq s \right\}; \\ \rho_{s}^{-} &= \inf \left\{ \mathbf{v}^{\top} \nabla^{2} f\left(\mathbf{\Sigma}\right) \mathbf{v} \mid \left\|\mathbf{v}\right\|_{2}^{2} = 1, \left\|\mathbf{v}\right\|_{0} \leq s \right\}. \end{split}$$

Moreover, we define $\kappa_s = \rho_s^+/\rho_s^-$ as the s-sparse condition number.

The sparse eigenvalue condition is widely studied in the high-dimensional statistics. And it is closely related to the restricted strongly convex, restricted smoothness and restricted eigenvalues properties [32]. Such condition have been employed by [33]-[35].

Definition 5. Define a local region of Σ^* by

$$\mathcal{B}(\mathbf{\Sigma}^*, r) = \{\mathbf{\Sigma} \succ 0 \mid ||\mathbf{\Sigma} - \mathbf{\Sigma}^*||_{\mathsf{F}} \le r\}.$$

In the following, we will show that $f(\Sigma)$ satisfies local strong convexity and local strong smoothness over a sparse domain. And in our analysis, we set the radius r as $\frac{\rho_{2s^*+2\tilde{s}}}{4\tau\kappa}$.

Assumption 6. Given $\Sigma \in \mathcal{B}(\Sigma^*, r)$ for a generic constant r, there exists a generic constant C_1 such that $\nabla^2 f(\Sigma)$ satisfies the sparse eigenvalue properties with parameters $\rho_{2s^*+2\widetilde{s}}^-$ and $\rho_{2s^*+2\widetilde{s}}^+$ satisfying

$$0 < \rho_{2s^*+2\widetilde{s}}^- < \rho_{2s^*+2\widetilde{s}}^+ < +\infty$$
 with $\widetilde{s} \ge C_1 \kappa_{2s^*+2\widetilde{s}}^2 s^*$ and $\kappa_{2s^*+2\widetilde{s}} = \rho_{2s^*+2\widetilde{s}}^+ / \rho_{2s^*+2\widetilde{s}}^-$

Assumption 6 shows that $\nabla^2 f(\Sigma)$ has bounded largest and non-zero smallest sparse eigenvalues, under the condition that Σ is adequately sparse and proximate to Σ^* . This employment of similar conditions has become a common practice in tackling high-dimensional problems [36]–[39]. Under our design, we confirm the existence of sparse eigenvalue. We refer readers to [40]–[46] for further details.

Assumption 7 (Local Restricted Hessian Smoothness). *Referencing* \widetilde{s} *introduced in Assumption 6. Constant* $L_{2s^*+2\widetilde{s}}$ *and* r *exist such that for any* $\Sigma, \Sigma' \in \mathcal{B}(\Sigma^*, r)$ *with* $\|\Sigma_{\overline{\mathcal{S}}}\|_0 \leq \widetilde{s}$ *and* $\|\Sigma'_{\overline{\mathcal{S}}}\|_0 \leq \widetilde{s}$, *the following inequality holds:*

$$\sup_{\mathbf{u} \in \Delta, \|\mathbf{u}\|_{\mathsf{F}} = 1} \|\mathbf{u}\|_{\nabla^2 f(\mathbf{\Sigma}') - \nabla^2 f(\mathbf{\Sigma})}^2 \le L_{2s^* + 2\widetilde{s}} \left\| \mathbf{\Sigma} - \mathbf{\Sigma}' \right\|_{\mathsf{F}}^2,$$

where $\Delta = \{\mathbf{u} \mid \mathsf{supp}(\mathbf{u}) \subseteq (\mathsf{supp}(\mathbf{\Sigma}) \cup \mathsf{supp}(\mathbf{\Sigma}'))\}$ and

$$\begin{split} & \left\| \mathbf{u} \right\|_{\nabla^{2} f(\mathbf{\Sigma}') - \nabla^{2} f(\mathbf{\Sigma})}^{2} \\ = & \mathsf{vec}^{\top} \left(\mathbf{u} \right) \left(\nabla^{2} f\left(\mathbf{\Sigma}'\right) - \nabla^{2} f\left(\mathbf{\Sigma}\right) \right) \mathsf{vec} \left(\mathbf{u} \right). \end{split}$$

Assumption 7 ensures that $\nabla^2 f(\Sigma)$ exhibits Lipschitz continuity in the vicinity of Σ^* across a sparse domain.

In the following, we specify some mild conditions for the true covariance matrix. We represent the true covariance matrix as Σ^* , and identify the support set of Σ^* as $\mathcal{S}^* = \{(j,k) \mid \Sigma_{jk}^* \neq 0\}$, with s^* denoting its size, that is, $s^* = |\mathcal{S}^*|$.

Assumption 8. Given the true covariance matrix Σ^* , there exist generic constants α , c such that

$$\left\| \mathbf{\Sigma}_{\mathcal{S}^*}^* \right\|_{\min} = \min_{(j,k) \in \mathcal{S}^*} \left| \mathbf{\Sigma}_{jk}^* \right| \ge (\alpha + c) \, \lambda \gtrsim \lambda,$$

where α and c being constants introduced in Assumption (2).

Assumption 8 guarantees the minimum signal strength condition, which is frequently utilized in the study of non-convex penalized regression problems [23], [24], [47]. In our design case, the tuning parameter λ can be taken to be the order of $\sqrt{\frac{\log d}{m}}$ that could not be too large when the measurements m increases, which guarantees that the estimators are close enough to the true model parameter. Taking the signal strength into account, we can obtain the oracle rate of convergence.

Assumption 9. For the true covariance matrix Σ^* , there exists some constant $\kappa \geq 1$ such that

$$0 < \frac{1}{\kappa} \le \lambda_{\mathsf{min}}\left(\boldsymbol{\Sigma}^*\right) \le \lambda_{\mathsf{max}}\left(\boldsymbol{\Sigma}^*\right) \le \kappa < \infty.$$

The condition regarding the smallest and largest eigenvalues of the true covariance matrix in Assumption 9 is standard within the existing research on the sparse covariance matrix estimation problem [48].

Assumption 10. At every stage of addressing the convex relaxed subproblem for all $q \ge 1$, we set

$$\varepsilon = \frac{C_3}{\sqrt{m}} \le \frac{\lambda}{8},$$

where C_3 is a predetermined small constant.

Assumption 10 ensures that the solution $\widetilde{\Sigma}^{(q)}$ obtained at each stage, for all $q \geq 1$, achieves adequate precision. This level of accuracy is essential for the convergence analysis of multistage convex relaxation.

B. Statistical Guarantees and Consequences

Now we introduce the primary result, illustrating the contraction property of the solution path $\left\{\widetilde{\Sigma}^{(q)}\right\}_{q\geq 1}$. For simplicity, we denote the following symbols. For functionals f(n) and g(n), $f(n)\gtrsim g(n)$ if $f(n)\geq cg(n)$, $f(n)\lesssim g(n)$ if $f(n)\leq Cg(n)$, and $f(n)\asymp g(n)$ if $cg(n)\leq f(n)\leq Cg(n)$ for some constant c and c. c0 is used to denote bounded in probability.

Theorem 11 (Contraction Property). Under the Assumptions 1, 6, 8 and 9, the ε -optimal solution $\widetilde{\Sigma}^{(q)}$ adheres to the following bound:

$$\begin{split} \left\| \widetilde{\boldsymbol{\Sigma}}^{(q)} - \boldsymbol{\Sigma}^* \right\|_{\mathsf{F}} &\leq \frac{1}{\rho_{2s^* + 2\widetilde{s}}^{-}} \underbrace{\| (\nabla f \left(\boldsymbol{\Sigma}^* \right))_{\mathcal{S}^*} \|_{\mathsf{F}}}_{oracle \ rate} \\ &+ \frac{1}{\rho_{2s^* + 2\widetilde{s}}^{-}} \underbrace{\varepsilon \sqrt{s^*}}_{optimization \ error} + \underbrace{\delta}_{contraction} \underbrace{\| \widetilde{\boldsymbol{\Sigma}}^{(q-1)} - \boldsymbol{\Sigma}^* \|_{\mathsf{F}}}_{contraction} \end{split}$$

for $1 \le q \le Q$, where $\delta \in (0,1)$ is the contraction factor.

Remark 12. The oracle rate describes the statistical convergence rate for an oracle estimator, which has prior knowledge the true support set \mathcal{S}^* . The oracle estimator $\widehat{\Sigma}^O$ is defined as

$$\widehat{\boldsymbol{\Sigma}}^{O} = \arg\min_{\boldsymbol{\Sigma}: \boldsymbol{\Sigma}_{\overline{S}^*} = \boldsymbol{0}} f(\boldsymbol{\Sigma}).$$

According to the definition, it follows quite straightforwardly that $\widehat{\Sigma}^O$ fulfills the condition $\left\|\widehat{\Sigma}^O - \Sigma^*\right\|_{\mathsf{F}} \lesssim \left\|\left(\nabla f\left(\Sigma^*\right)\right)_{\mathcal{S}^*}\right\|_{\mathsf{F}}$, indicating a close approximation to Σ^* within the confines of the true support set's gradient norms.

Theorem 11 reveals the estimation discrepancy between the ε -optimal solution $\widetilde{\Sigma}^{(q)}$ and the actual parameter Σ^* is bounded by three components: the oracle rate, the optimization error, and a contraction term. Following this, we detail the precise statistical convergence rate applicable within our framework.

Corollary 13. Suppose that Assumptions 1, 6, 8, 9 and 10 hold. If $\lambda \asymp \sqrt{\frac{\log d}{m}}$, $\tau \lesssim \sqrt{\frac{1}{m}} \left\| (\boldsymbol{\Sigma}^*)^{-1} \right\|_{\max}^{-1}$, the ε -optimal solution $\widetilde{\boldsymbol{\Sigma}}^{(1)}$ satisfies

$$\left\|\widetilde{\boldsymbol{\Sigma}}^{(1)} - \boldsymbol{\Sigma}^*\right\|_{\mathsf{F}} \lesssim \sqrt{\frac{s^* \log d}{m}}$$

with high probability.

Corollary 13 naturally follows from Theorem 11 when q=1. Furthermore, the contraction property, as discussed in Theorem 11, is induced by the MM-based multistage convex relaxation algorithm. To attain the oracle rate, the optimization error must be strategically selected so that $\varepsilon \leq \frac{\left\|(\nabla f(\mathbf{\Sigma}^*))_{\mathcal{S}^*}\right\|_{\mathsf{F}}}{\sqrt{s^*}}$ and the parameter Q should be sufficiently large. This leads us to the subsequent conclusion.

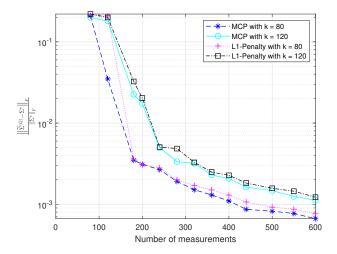


Fig. 1. The mean relative error of the estimated covariance matrices for different sparsity level with noise.

Corollary 14. Suppose that Assumptions 1, 6, 7, 8, 9, and 10 hold. If $\lambda \asymp \sqrt{\frac{\log d}{m}}$, $\tau \lesssim \sqrt{\frac{1}{m}} \left\| (\mathbf{\Sigma}^*)^{-1} \right\|_{\max}^{-1}$, and $Q \gtrsim \log (\lambda \sqrt{m}) \gtrsim \log \log d$, then the ε -optimal solution $\widetilde{\mathbf{\Sigma}}^{(Q)}$ satisfies

$$\left\|\widetilde{\boldsymbol{\Sigma}}^{(Q)} - \boldsymbol{\Sigma}^*\right\|_{\mathsf{F}} = \mathcal{O}_p\left(\sqrt{\frac{s^*}{m}}\right).$$

Corollary 14 is a direct consequence of Theorem 11, which suggests that under minimal assumptions, solving no more than approximately $\log \log d$ convex problems is sufficient to reach the oracle rate $\sqrt{\frac{s^*}{m}}$.

V. NUMERICAL EXPERIMENTS

We compare our proposal with the existing estimator, focusing specifically on the estimator that incorporates an ℓ_1 penalty. We conduct a series of Monte Carlo simulations utilizing matrices of dimensions 50×50 . The tuning parameters λ and τ were selected via five-fold cross-validation. Initially, positive symmetric sparse covariance matrices are generated utilizing the "sprandsym" function in MATLAB, defined by a sparsity value k, which represents the quantity of non-zero elements within the matrix. The non-convex penalty function chosen for these experiments is the MCP, with a constant setting of b=2 across all trials.

For each scenario characterized by a specific (m, \mathbf{k}) pair, we execute 10 independent iterations. The outcomes are depicted in Figure 1, which illustrates the mean relative error of the estimated covariance matrices as a function of m across varying levels of sparsity. Each measurement includes additive noise, achieved by generating η_i from a normal distribution $\sigma \cdot \mathcal{N} (0,1)$ with $\sigma = 10^{-4}$. From the simulation results, it is evident that the estimator with MCP performs better than the ℓ_1 penalty-based estimator in terms of Frobenius norm error. This supports our theoretical findings that non-convex penalty can reduce the covariance estimation error.

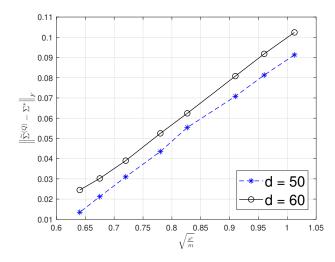


Fig. 2. The oracle rate for different measurements and different dimension of Σ^* .

Fig. 2 illustrates the performance of oracle rate in estimating a covariance matrix from quadratic measurement model. The y-axis represents the deviation between an estimated covariance matrix and the true covariance matrix, measured using the Frobenius norm. The x-axis is a scaled measure related to the support size of true covariance matrix and the measurement complexity, indicated by $\sqrt{\frac{s^*}{m}}$. Two lines are plotted to show the estimation error for two different dimensions $(d = \{50, 60\})$ under the sparsity level of k = 80. As the measurement size increase, the error between the estimated and true covariance matrix decreases. The black line (d = 60) shows a slightly higher error trend than the blue line (d = 50), suggesting that higher-dimensional covariance tends to have larger deviations from the true matrix. This visual representation provides insight into how the quality of covariance matrix estimation varies with the oracle rate.

VI. CONCLUSION AND DISCUSSION

This paper introduces a new sample strategy named covariance matrix sensing, including sensing and compressing signals and estimating the covariance matrix from a small number of measurements. Specifically, we estimate large sparse covariance matrices from quadratic samplings through a non-convex penalty, detailing both its theoretical results and empirical validations. We demonstrate that our estimators achieve superior statistical convergence rates when contrasted with current methodologies.

And all theoretical derivations and the suggested algorithm can be seamlessly applied to the bilinear measurement model where $y_i = a_i^{\top} \Sigma b_i$, with a_i and b_i being independently produced sensing vectors. This demonstrates that our outcomes are equally applicable to this asymmetric sensing model too.

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