SOLVING LOW-RANK SEMIDEFINITE PROGRAMS VIA MANIFOLD OPTIMIZATION *

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Abstract. We propose a manifold optimization approach to solve linear semidefinite programs (SDP) with low-rank solutions. This approach incorporates the augmented Lagrangian method and the Burer-Monteiro factorization, and features the adaptive strategies for updating the factorization size and the penalty parameter. We prove that the present algorithm can solve SDPs to global optimality, despite of the non-convexity brought by the Burer-Monteiro factorization. Along with the algorithm, we release an open-source SDP solver ManiSDP. Comprehensive numerical experiments demonstrate that ManiSDP achieves state-of-the-art in terms of efficiency, accuracy and scalability, and is faster than several advanced SDP solvers (MOSEK, SDPLR, SDPNAL+, STRIDE) by up to orders of magnitudes on a variety of SDPs. The largest SDP solved by ManiSDP (in about 8.5 hours with maximal KKT residue 3.5e-13) is the second-order moment relaxation of a binary quadratic program with 120 variables, which has matrix dimension 7261 and contains 17, 869, 161 affine constraints.

Key words. semidefinite programming, low-rank solution, polynomial optimization, moment relaxation, Burer-Monteiro factorization, augmented Lagrangian method, manifold optimization

AMS subject classifications. Primary, 90C22; Secondary, 90C23,90C30

1. Introduction. In this paper, we aim to efficiently solve the following semi-definite programming (SDP) problem:

$$\begin{cases} \inf_{X\succeq 0} & \langle C,X\rangle\\ \text{s.t.} & \mathcal{A}(X)=b, \mathcal{B}(X)=d, \end{cases}$$

where $\mathcal{A}: \mathbb{S}_n \to \mathbb{R}^m$, $\mathcal{B}: \mathbb{S}_n \to \mathbb{R}^l$ are linear maps, $b \in \mathbb{R}^m$, $d \in \mathbb{R}^l$. In (SDP) the linear constraints $\mathcal{A}(X) = b$ are arbitrary while the linear constraints $\mathcal{B}(X) = d$ are assumed to impose certain manifold structures on the domain if present. Moreover, we assume that (SDP) admits a low-rank solution X^* , i.e., $\operatorname{rank}(X^*) \ll n$.

Such (SDP) arises from a diverse set of fields, e.g., systems and control [25], signal processing [15, 23], optimal power flow [4], matrix completion [10], quantum steering [12], computer vision [38], to just name a few; see also the surveys [30, 36] and references therein. It also serves as a tractable convex relaxation for many difficult (usually NP-hard) non-convex optimization problems, e.g., quadratically constrained quadratic programs [23], combinatorial optimization problems [17], polynomial optimization problems [22].

In view of the wide applications, a large amount of efforts have been dedicated to solving (SDP) over the past decades and lots of algorithms have been developed from various angles. For small/medium-scale SDPs, interior-point methods are believed to be the most accurate, efficient and robust algorithms [3, 29]. However, for large-scale SDPs interior-point methods are no longer reliable because of their extensive memory occupation and high computational costs in solving a large and dense linear system at each iteration. Aiming to tackle large-scale SDPs, Yang, Sun and Toh proposed

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an augmented Lagrangian algorithm combined with the semismooth Newton method. The related solver SDPNAL+ has been shown to have good scalability on *degenerate* SDPs [41]. There are also quite a few attempts to overcome the memory issue by relying on first-order methods. For instance, Wen et al. proposed an alternating-direction augmented Lagrangian method (ALM) [35]. The main drawback of first-order methods is that they can hardly achieve high accuracy.

In many cases, large-scale SDPs possess certain structures, e.g., chordal sparsity, constant trace, admitting low-rank solutions. Such structures can be exploited to speed up computation either in interior-point methods or in first-order methods. Zhang and Lavaei proposed to exploit chordal sparsity for interior-point methods via dualized clique tree conversion [44]. Zheng et al. [45] and Garstka et al. [16] proposed to exploit chordal sparsity in a framework of the alternating direction methods of multipliers (ADMM). By exploiting the constant trace property, Helmberg and Rendl proposed a spectral bundle method for solving (SDP) [19]. More recently, Yurtsever et al. proposed to exploit the constant trace property in a conditional-gradient-based augmented Lagrangian framework combined with a matrix sketching technique [42]. The property of admitting low-rank solutions can be exploited in the framework of interior-point methods [5, 18, 43] or in the framework of operator splitting methods [27]. For SDPs arising as convex relaxations of polynomial optimization problems and admitting rank-one solutions, Yang et al. proposed a projected gradient method accelerated by local search [39], which is scalable on a variety of SDPs emerging from computer vision problems [38].

Another notable approach taking the low-rank property into account is to perform a factorization $X = YY^{\mathsf{T}}$ with $Y \in \mathbb{R}^{n \times p}, p \ll n$ so that (SDP) becomes a non-convex QCQP (called the factorized problem), which is called the Burer-Monteiro factorization in the literature [8, 9], leading to a low-rank SDP solver SDPLR. For the factorized problem, nonlinear programming tools can find a locally optimal solution at much less cost, and under mild conditions, this local solution is guaranteed to be globally optimal whenever $\frac{p(p+1)}{2} > m$ [7]. Hence for the approach being efficient, m is required not to be too large. Following Burer and Monteiro's idea, Journée et al. reformulated (SDP) as a nonlinear program on a manifold under the assumption that m=0 (no arbitrary linear constraint) and the feasible set $\mathcal{N}_p = \{Y \in \mathbb{R}^{n \times p} \mid \mathcal{B}(YY^{\mathsf{T}}) = d\}$ is a smooth manifold [20]. Consequently, the nonlinear program becomes a Riemannian optimization problem which can be efficiently solved by off-the-shelf Riemannian optimization tools. They also provided a strategy for escaping from saddle points so that global convergence can be guaranteed. More recently, this approach has been extended to include arbitrary linear constraints A(X) = b and to handle nonsmooth objective functions within an augmented Lagrangian framework in which the Riemannian semismooth Newton method is employed to solve the augmented Lagrangian subproblem [34]. See also [28] for the extension of the approach to graph equipartition SDPs and [26] for the extension to the regularized convex matrix optimization problem.

Contributions. Our contributions are threefold:

(1) We present a manifold optimization algorithm for solving (SDP) by adopting the Burer-Monteiro factorization and the idea of Journée et al.. More concretely, our algorithm employs an ALM framework to handle arbitrary affine constraints and applies the Burer-Monteiro factorization to the ALM subproblem in order to exploit the low-rank property. The ALM subproblem is then recast as a manifold optimization problem which is solved using the Riemannian trust-region method. To circumvent

the non-convexity introduced by the Burer-Monteiro factorization, we design a strategy (inspired by Journée et al.) for escaping from saddle points. The overall algorithm possesses two distinguished features: (i) the factorization size p is dynamically adjusted so that the number of decision variables of the ALM subproblem is as small as possible; (ii) a novel strategy of self-adaptively updating the penalty parameter which allows to decrease the penalty parameter is embedded. These two features significantly improve the practical performance of the algorithm.

- (2) Despite of the non-convexity, we prove global convergence of the proposed algorithm, that is, we show that under certain conditions, the output of the algorithm satisfies the KKT conditions of (SDP) up to a small tolerance.
- (3) We release an open-source SDP solver named ManiSDP that implements the proposed algorithm. Extensive numerical experiments were performed to benchmark the solver, which demonstrate that ManiSDP is accurate, efficient and scalable on a variety of SDPs with low-rank solutions, and (substantially) outperforms a few popular SDP solvers.

The rest of the paper is organized as follows. In Section 2, we collect notations and some preliminary results. In Section 3, we give the augmented Lagrangian framework with the Burer-Monteiro factorization. In Section 4, we describe some computational details on solving the ALM subproblem with the Riemannian trust-region method. In Section 5, we describe the strategies of updating the factorization size and the penalty parameter, and prove global convergence of the algorithm. Results of numerical experiments are provided in Section 6. Conclusions are made in Section 7.

2. Notation and preliminaries. \mathbb{R} (resp. \mathbb{N}) denotes the set of real numbers (resp. nonnegative integers). For a positive integer n, let $[n] := \{1, 2, \dots, n\}$. Let \mathbb{S}_n (resp. \mathbb{S}_n^+) denote the set of (resp. positive semidefinite/PSD) symmetric matrices of size n. We use $\mathrm{Tr}(A)$ (resp. A^{T}) to denote the trace (resp. transpose) of a matrix $A \in \mathbb{R}^{n \times m}$. For two matrix $A, B \in \mathbb{R}^{n \times m}$, the inner product is defined as $\langle A, B \rangle = \mathrm{Tr}(A^{\mathsf{T}}B)$. For $A \in \mathbb{R}^{n \times n}$, diag(A) denotes the diagonal of A and Diag(A) denotes the diagonal matrix with the same diagonal as A. For a vector v, ||v|| is the 2-norm of v and for a matrix A, $||A||_F$ is the Frobenius norm of A. For a function f(x), we write $\nabla f(x)$ (resp. grad f(x)) for the Euclidean (resp. Riemannian) gradient, and write $\nabla^2 f(x)[u]$ (resp. Hess f(x)[u]) for the Euclidean (resp. Riemannian) Hessian acting on u. For a set E, |E| denotes its cardinality.

Let us consider (SDP) with $\mathcal{A}(X) := (\langle A_i, X \rangle)_{i=1}^m$ and $\mathcal{B}(X) := (\langle B_i, X \rangle)_{i=1}^l$ for $A_i, B_i \in \mathbb{S}_n$. Let $\mathcal{A}^* : \mathbb{R}^m \to \mathbb{S}_n$ be the adjoint operator of \mathcal{A} defined as $\mathcal{A}^*(y) := \sum_{i=1}^m y_i A_i$ for $y \in \mathbb{R}^m$; similarly, let $\mathcal{B}^* : \mathbb{R}^l \to \mathbb{S}_n$ be the adjoint operator of \mathcal{B} defined as $\mathcal{B}^*(z) := \sum_{i=1}^l z_i B_i$ for $z \in \mathbb{R}^l$. Throughout the paper, we assume that strong duality holds for (SDP).

LEMMA 2.1. A matrix $X \in \mathbb{S}_n$ is a minimizer of (SDP) if and only if there exist Lagrange multipliers $y \in \mathbb{R}^m$ and $z \in \mathbb{R}^l$ such that

$$\mathcal{A}(X) = b,$$

$$\mathcal{B}(X) = d,$$

$$(2.1c) X \succeq 0,$$

(2.1d)
$$S := C - \mathcal{A}^*(y) - \mathcal{B}^*(z) \succeq 0,$$

$$(2.1e) XS = 0.$$

Proof. These are the standard KKT conditions for (SDP).

- 3. An augmented Lagrangian framework with the Burer-Monteiro factorization. In this section, we introduce an augmented Lagrangian framework combined with the Burer-Monteiro factorization for solving (SDP).
- 3.1. An augmented Lagrangian framework for solving semidefinite programs. Let us denote

(3.1)
$$\mathcal{M} := \{ X \in \mathbb{R}^{n \times n} \mid \mathcal{B}(X) = d, X \succeq 0 \}.$$

Note that if l = 0 (i.e., the constraints $\mathcal{B}(X) = d$ are not present), then $\mathcal{M} = \mathbb{S}_n^+$. Other typical choices of \mathcal{M} are

(Unit-trace)
$$\mathcal{M} = \{ X \in \mathbb{R}^{n \times n} \mid X \succeq 0, \text{Tr}(X) = 1 \},$$
(Unit-diagonal)
$$\mathcal{M} = \{ X \in \mathbb{R}^{n \times n} \mid X \succeq 0, \text{diag}(X) = \mathbf{1} \}.$$

Remark 3.1. For the case that X has a constant trace c, we can scale X by the factor $\frac{1}{c}$ to match the (Unit-trace) case.

Then (SDP) can be equivalently written as

(SDP-M)
$$\begin{cases} \inf_{X \in \mathcal{M}} & \langle C, X \rangle \\ \text{s.t.} & \mathcal{A}(X) = b. \end{cases}$$

We denote by \mathcal{N}_p the set of matrices after applying the Burer-Monteiro factorization to \mathcal{M} , i.e.,

$$(3.2) \mathcal{N}_p := \{ Y \in \mathbb{R}^{n \times p} \mid YY^{\mathsf{T}} \in \mathcal{M} \} = \{ Y \in \mathbb{R}^{n \times p} \mid \mathcal{B}(YY^{\mathsf{T}}) = d \}.$$

For $\mathcal{M} = \mathbb{S}_n^+$ or \mathcal{M} in (Unit-trace), (Unit-diagonal), the corresponding \mathcal{N}_p are

(Euclidean)
$$\mathcal{N}_p = \{Y \in \mathbb{R}^{n \times p}\},$$
(Sphere)
$$\mathcal{N}_p = \{Y \in \mathbb{R}^{n \times p} \mid ||Y||_F = 1\},$$
(Oblique)
$$\mathcal{N}_p = \{Y \in \mathbb{R}^{n \times p} \mid ||Y(i,:)|| = 1, i = 1, \dots, n\},$$

where Y(i,:) stands for the *i*-th row of Y.

We further make the following assumptions on (SDP-M):

Assumption 3.2. (SDP-M) admits a low-rank optimal solution.

Assumption 3.3. \mathcal{N}_p is a submanifold embedded in the Euclidean space.

Assumption 3.4. Either l=0,1, or the matrices $\{B_i\}_{i=1}^l$ satisfy $B_iB_j=0$ for any $1 \le i \ne j \le l$.

Assumption 3.5. $d_i \neq 0$ for i = 1, ..., l.

It is clear that for \mathcal{N}_p in (Euclidean) or (Sphere) with $p \geq 1$, or \mathcal{N}_p in (Oblique) with $p \geq 2$, Assumptions 3.3–3.5 are satisfied.

(SDP-M) can be solved with the ALM as presented in Algorithm 3.1.

The augmented Lagrangian function associated with (SDP-M) is defined by

(3.4)
$$L_{\sigma}(X,y) = \langle C, X \rangle - y^{\mathsf{T}}(\mathcal{A}(X) - b) + \frac{\sigma}{2} \|\mathcal{A}(X) - b\|^{2}.$$

Algorithm 3.1

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Input: A, b, B, \overline{d}, C, \sigma_0 > 0
Output: (X, y, S)
 1: k \leftarrow 0, y^0 \leftarrow 0;
 2: while stopping criteria do not fulfill do
        Solve the ALM subproblem (3.5) to obtain X^{k+1};
        y^{k+1} \leftarrow y^k - \sigma_k(\mathcal{A}(X^{k+1}) - b);
 4:
        Compute \sigma_{k+1};
 5:
        k \leftarrow k + 1:
 7: end while
 8: return (X^k, y^k, S^k);
```

At the k-th iteration of the ALM, we need to solve the subproblem

(3.5)
$$\min_{X \in \mathcal{M}} \Phi_k(X) := L_{\sigma^k}(X, y^k).$$

The following lemma characterizes the optimality conditions of (3.5).

LEMMA 3.6. A matrix $X \in \mathcal{M}$ is a minimizer of (3.5) if and only if there exist Lagrange multipliers $z \in \mathbb{R}^l$ such that

(3.6a)
$$S := \nabla \Phi_k(X) - \mathcal{B}^*(z) \succeq 0,$$

$$(3.6b) XS = 0,$$

where $\nabla \Phi_k(X) = C + \sigma_k \mathcal{A}^*(\mathcal{A}(X) - b - y^k/\sigma_k)$. Furthermore, under Assumptions 3.4-3.5, the vector z is unique and has a closed-form expression:

(3.7)
$$z_i := \frac{\operatorname{Tr}(B_i X \nabla \Phi_k(X))}{\operatorname{Tr}(B_i^2 X)} \text{ for } i = 1, \dots, l,$$

Proof. As (3.5) is convex, X is a minimizer if and only if the KKT conditions hold, i.e., there exist $z \in \mathbb{R}^l$ such that

$$(3.8a) \mathcal{B}(X) = d,$$

$$(3.8b) X \succeq 0,$$

$$(3.8c) S \succeq 0,$$

$$(3.8d) XS = 0.$$

Since $X \in \mathcal{M}$, (3.8a) and (3.8b) are valid by definition, and so X is a minimizer if and only if (3.6a) and (3.6b) hold for some $z \in \mathbb{R}^l$. To get the closed-form expression of z, we note that XS = 0 can be rewritten as

$$(3.9) X\nabla\Phi_k(X) - \sum_{j=1}^l XB_j z_j = 0.$$

For each $i \in [l]$, multiplying (3.9) by B_i and noting $B_iB_j = 0$ for $i \neq j$, we get

$$(3.10) X\nabla\Phi_k(X)B_i - XB_i^2 z_i = 0.$$

Let $X = YY^{\mathsf{T}}$ for some $Y \in \mathcal{N}_p$. We have

$$\operatorname{Tr}(XB_{i}^{2}) = \operatorname{Tr}(Y^{\mathsf{T}}B_{i}^{2}Y) = \|B_{i}Y\|_{F}^{2} \neq 0,$$

where the last inequality is because B_iY cannot be a zero matrix since $d_i \neq 0$. Then (3.7) follows by taking trace of (3.10).

Remark 3.7. For \mathcal{M} in (Unit-trace), (3.7) becomes

$$(3.11) z = \operatorname{Tr}(X \nabla \Phi_k(X)),$$

and for \mathcal{M} in (Unit-diagonal), (3.7) becomes

(3.12)
$$z = \operatorname{diag}(X \nabla \Phi_k(X)).$$

COROLLARY 3.8. Given a matrix $X \in \mathcal{M}$, let $S := \nabla \Phi_k(X) - \mathcal{B}^*(z)$ with z being given by (3.7). Then X is a minimizer of the convex subproblem (3.5) if and only if $S \succeq 0$.

Proof. This is an immediate consequence of Lemma 3.6.

3.2. The Burer-Monteiro factorization. To exploit the fact that (SDP-M) admits a low-rank optimal solution (Assumption 3.2), we now apply the Burer-Monteiro factorization to $\Phi_k(X)$ and define $\Psi_k(Y) = \Phi_k(YY^{\mathsf{T}})$ for $Y \in \mathbb{R}^{n \times p}$. Consequently, the convex subproblem (3.5) becomes the nonconvex factorized subproblem:

$$(3.13) \qquad \min_{Y \in \mathcal{N}_p} \Psi_k(Y) = \langle C, YY^\mathsf{T} \rangle - (y^k)^\mathsf{T} (\mathcal{A}(YY^\mathsf{T}) - b) + \frac{\sigma_k}{2} \|\mathcal{A}(YY^\mathsf{T}) - b\|^2.$$

Then, we solve the nonconvex factorized subproblem (3.13) on the manifold \mathcal{N}_p instead of solving the convex subproblem (3.5) on \mathcal{M} . The global optimality condition of (3.13) can be characterized in terms of positive semidefiniteness of the matrix S as well.

PROPOSITION 3.9. Let $Y \in \mathcal{N}_p$, $X = YY^{\mathsf{T}} \in \mathcal{M}$, and z be given by (3.7). Let $S = \nabla \Phi_k(X) - \mathcal{B}^*(z)$. Then Y is a global minimizer of the nonconvex factorized subproblem (3.13) if and only if $S \succeq 0$.

Proof. Note that Y is a global minimizer of (3.13) if and only if X is a minimizer of the convex subproblem (3.5). Then the conclusion follows from Corollary 3.8.

THEOREM 3.10. A matrix $Y \in \mathcal{N}_p$ provides a minimizer $X = YY^{\mathsf{T}}$ of (SDP-M) if and only if $\mathcal{A}(X) = b$ and $S = \nabla \Phi_k(X) - \mathcal{B}^*(z) \succeq 0$ with z being given by (3.7).

Proof. If X is a minimizer of (SDP-M), then $\mathcal{A}(X) = b$ and X is a minimizer of (3.5). So by Corollary 3.8, $S \succeq 0$. Conversely, by Corollary 3.8, $S \succeq 0$ implies X is a minimizer of (3.5) and so XS = 0. Moreover, from $\mathcal{A}(X) = b$, we get $S = C - \mathcal{A}^*(y^k) - \mathcal{B}^*(z)$. Then we see that (X, y^k, z, S) satisfies the KKT conditions (2.1), which yields that X is a minimizer of (SDP-M).

- 4. Solve the ALM subproblem by the Riemannian trust-region method. Since \mathcal{N}_p is assumed to be a manifold, we can solve the nonconvex factorized subproblem (3.13) by off-the-shelf efficient manifold optimization methods. Here we choose the Riemannian trust-region method for its superior performance in both accuracy and efficiency, though other choices are also possible.
- **4.1. The Riemannian trust-region method.** In this subsection, we calculate the ingredients that are necessary to perform optimization on the manifold \mathcal{N}_p via the Riemannian trust-region method. For a detailed introduction to the Riemannian trust-region method, we refer the reader to [1]. First of all, we note that at a point $Y \in \mathcal{N}_p$, the tangent space of \mathcal{N}_p is given by

$$(4.1) T_Y \mathcal{N}_p = \{ U \in \mathbb{R}^{n \times p} \mid \mathcal{B}(UY^{\mathsf{T}}) = 0 \},$$

and the normal space to \mathcal{N}_p is given by

$$(4.2) N_Y \mathcal{N}_p = \left\{ \mathcal{B}^*(u) Y = \sum_{i=1}^l u_i B_i Y \middle| u \in \mathbb{R}^l \right\}.$$

LEMMA 4.1. Let Y be a point on \mathcal{N}_p . The orthogonal projector $P_Y: \mathbb{R}^{n \times p} \to T_Y \mathcal{N}_p$ is given by

$$(4.3) P_Y(U) = U - \mathcal{B}^*(u)Y = U - \sum_{i=1}^l u_i B_i Y, \text{ for } U \in \mathbb{R}^{n \times p},$$

where $u \in \mathbb{R}^l$ is determined by

(4.4)
$$u_i = \frac{\operatorname{Tr}(B_i U Y^{\mathsf{T}})}{\operatorname{Tr}(B_i^2 Y Y^{\mathsf{T}})}, \text{ for } i = 1, \dots, l.$$

Proof. Any matrix $U \in \mathbb{R}^{n \times p}$ admits a unique decomposition $U = U_{T_Y \mathcal{N}_p} + U_{N_Y \mathcal{N}_p}$, where $U_{\mathcal{X}}$ belongs to the space \mathcal{X} . So the projection $P_Y(U)$ can be assumed to be of form $P_Y(U) = U - \mathcal{B}^*(u)Y$ for some $u \in \mathbb{R}^l$. Since $P_Y(U) \in T_Y \mathcal{N}_p$, it holds

$$\mathcal{B}(P_Y(U)Y^{\mathsf{T}}) = \mathcal{B}(UY^{\mathsf{T}}) - \mathcal{B}(\mathcal{B}^*(u)YY^{\mathsf{T}}) = 0,$$

i.e., $\operatorname{Tr}(B_iUY^{\mathsf{T}}) = \operatorname{Tr}(B_i\mathcal{B}^*(u)YY^{\mathsf{T}})$ for $i = 1, \ldots, l$. This yields (4.4) as $B_iB_j = 0$ for $i \neq j$ by Assumption 3.4.

PROPOSITION 4.2. Consider the nonconvex factorized subproblem (3.13). Let $X = YY^{\mathsf{T}}$, $\widetilde{S} = \nabla \Phi_k(X)$ and $S = \widetilde{S} - \mathcal{B}^*(z)$ with z being given in (3.7). Then the Riemannian gradient at Y is given by

(4.5)
$$\operatorname{grad}\Psi_k(Y) = 2SY.$$

For $U \in T_Y \mathcal{N}_p$, let $\widetilde{H} = \nabla^2 \Psi_k(Y)[U] = 2(\widetilde{S}U + \sigma_k \mathcal{A}^* (\mathcal{A}(YU^{\mathsf{T}} + UY^{\mathsf{T}}))Y)$. Then the Riemannian Hessian is given by

(4.6)
$$\operatorname{Hess} \Psi_k(Y)[U] = P_Y(\widetilde{H}) - 2\mathcal{B}^*(z)U + 2\mathcal{B}^*(u)Y,$$

where

(4.7)
$$u_i = \frac{\operatorname{Tr}(B_i^2 U Y^{\dagger}) z_i}{\operatorname{Tr}(B_i^2 X)}, \text{ for } i = 1, \dots, l.$$

In particular, for \mathcal{N}_p in (Euclidean), we have

(4.8)
$$\operatorname{Hess} \Psi_k(Y)[U] = \widetilde{H};$$

for \mathcal{N}_p in (Sphere), we have

(4.9)
$$\operatorname{Hess} \Psi_k(Y)[U] = \widetilde{H} - \operatorname{Tr}(\widetilde{H}Y^{\dagger})Y - 2\operatorname{Tr}(\widetilde{S}X)U;$$

for \mathcal{N}_p in (Oblique), we have

(4.10)
$$\operatorname{Hess} \Psi_k(Y)[U] = \widetilde{H} - \operatorname{Diag}(\widetilde{H}Y^{\mathsf{T}})Y - 2\operatorname{Diag}(\widetilde{S}X)U.$$

Proof. First note $\nabla \Psi_k(Y) = 2\widetilde{S}Y$. By (3) of [2], grad $\Psi_k(Y) = P_Y(\nabla \Psi_k(Y))$ and thus we can write

$$\operatorname{grad} \Psi_k(Y) = 2\widetilde{S}Y - 2\mathcal{B}^*(z)Y$$

for some $z \in \mathbb{R}^l$. As grad $\Psi_k(Y) \in T_Y \mathcal{N}_p$, we have

(4.11)
$$\mathcal{B}(\operatorname{grad}\Psi_k(Y)Y^{\mathsf{T}}) = 2\mathcal{B}((\widetilde{S} - \mathcal{B}^*(z))X) = 0.$$

Solving (4.11) for z we get exactly (3.7) and (4.5) then follows. By (10) of [2], it holds

(4.12)
$$\operatorname{Hess}\Psi_k(Y)[U] = P_Y(\nabla^2 \Psi_k(Y)[U]) + \mathfrak{A}_Y(U, P_Y^{\perp}(\nabla \Psi_k(Y))),$$

where \mathfrak{A}_Y is the Weingarten map at Y and $P_Y^{\perp} = I - P_Y$ is the orthogonal projector at Y to $N_Y \mathcal{N}_p$. Let $D_Y(\cdot)[U]$ be the directional derivative at Y along U. Then we have

$$\begin{split} \mathfrak{A}_Y(U,P_Y^\perp(\nabla\Psi_k(Y))) &= \mathfrak{A}_Y(U,2\mathcal{B}^*(z)Y) \\ &= -P_Y(D_Y(2\mathcal{B}^*(z)Y)[U]) \\ &= -2P_Y(\mathcal{B}^*(z)U) - 2P_Y(\mathcal{B}^*(D_Y(z)[U])Y) \\ &= -2\mathcal{B}^*(z)U + 2\mathcal{B}^*(u)Y, \end{split}$$

where we have used the fact that $P_Y(\mathcal{B}^*(z)U) = \mathcal{B}^*(z)U - \mathcal{B}^*(u)Y$ and $P_Y(\mathcal{B}^*(u')Y) = 0$ for any $u' \in \mathbb{R}^l$. (4.6) then follows.

The remaining conclusions of the proposition can be easily verified.

4.2. Escaping from saddle points. Since the subproblem (3.13) is nonconvex, in order to solve (3.13) to global optimality, it is then crucial to design a strategy for escaping from saddle points. We next show that we can always compute a second-order descent direction U to escape from saddle points whenever $S = \nabla \Phi_k(X) - \mathcal{B}^*(z) \not\succeq 0$.

LEMMA 4.3. For any $U \in \mathbb{R}^{n \times p}$ satisfying $YU^{\mathsf{T}} = 0$, it holds

$$(4.13) \langle U, \operatorname{Hess} \Psi_k(Y)[U] \rangle = 2 \operatorname{Tr}(U^{\mathsf{T}} S U).$$

Proof. The assumption $YU^{\intercal} = 0$ implies $U \in T_Y \mathcal{N}_p$. By (4.6) and Lemma 4.1, $\widetilde{H} = 2\widetilde{S}U$ with $\widetilde{S} = \nabla \Phi_k(X)$ and

$$\begin{aligned} \operatorname{Hess} \Psi_k(Y)[U] &= 2P_Y(\widetilde{S}U) - 2\mathcal{B}^*(z)U + 2\mathcal{B}^*(u)Y \\ &= 2\widetilde{S}U - 2\mathcal{B}^*(u')Y - 2\mathcal{B}^*(z)U + 2\mathcal{B}^*(u)Y \\ &= 2SU + 2\mathcal{B}^*(u - u')Y \end{aligned}$$

for z, u in Proposition 4.2 and some $u' \in \mathbb{R}^l$. Thus,

$$\langle U, \operatorname{Hess} \Psi_k(Y)[U] \rangle = 2\operatorname{Tr}(U^{\mathsf{T}}SU) + 2\operatorname{Tr}(U^{\mathsf{T}}\mathcal{B}^*(u-u')Y) = 2\operatorname{Tr}(U^{\mathsf{T}}SU).$$

THEOREM 4.4. Suppose $S = \nabla \Phi_k(X) - \mathcal{B}^*(z) \succeq 0$ with z being given by (3.7). Let $V \in \mathbb{R}^{n \times \delta}$ be a matrix whose columns consist of eigenvectors corresponding to negative eigenvalues of S. Then $U := [0_{n \times p}, V]$ is a second-order descent direction of (3.13) with $p := p + \delta$ at the point $Y := [Y, 0_{n \times \delta}]$, namely, U satisfies

$$(4.14) \langle U, \operatorname{grad} \Psi_k(Y) \rangle = 0, \quad \langle U, \operatorname{Hess} \Psi_k(Y)[U] \rangle < 0.$$

Proof. By construction, we have $YU^{\dagger} = 0$. Therefore, by (4.5),

$$\langle U, \operatorname{grad} \Psi_k(Y) \rangle = 2\operatorname{Tr}(U^{\mathsf{T}}SY) = 2\operatorname{Tr}(SYU^{\mathsf{T}}) = 0.$$

By Lemma 4.3,

$$\langle U, \operatorname{Hess} \Psi_k(Y)[U] \rangle = 2\operatorname{Tr}(U^{\mathsf{T}}SU) = 2\operatorname{Tr}(V^{\mathsf{T}}SV) < 0.$$

The following theorem tells us that by escaping from saddle points, we are able to find a global minimizer of the nonconvex factorized subproblem (3.13).

Theorem 4.5. A local minimizer $Y \in \mathcal{N}_p$ of the nonconvex factorized subproblem (3.13) provides a minimizer $X = YY^{\intercal}$ of the convex subproblem (3.5) if it is rank deficient, i.e., rank Y < p.

Proof. Let rank Y = r < p. We can write $Y = Y'P^{\intercal}$ for some full-rank matrices $Y' \in \mathbb{R}^{n \times r}$ and $P \in \mathbb{R}^{p \times r}$. Let $P_{\perp} \in \mathbb{R}^{p \times (p-r)}$ be the matrix whose columns form an orthogonal basis of the orthogonal complement of the column space of P such that $P^{\intercal}P_{\perp} = 0$. For any $V \in \mathbb{R}^{n \times (p-r)}$, let $U = VP_{\perp}^{\intercal}$. Then $UY^{\intercal} = VP_{\perp}^{\intercal}P(Y')^{\intercal} = 0$. As Y is a local minimizer of the nonconvex factorized subproblem (3.13), Hess $\Psi_k(Y)$ is positive semidefinite. Then by Lemma 4.3, we have

$$0 \le \langle U, \operatorname{Hess} \Psi_k(Y)[U] \rangle = 2 \operatorname{Tr}(U^{\mathsf{T}} S U) = 2 \operatorname{Tr}(V^{\mathsf{T}} S V).$$

The arbitrarity of V implies that S is positive semidefinite and so by Proposition 3.9, $Y \in \mathcal{N}_p$ is a global minimizer of (3.13), which further implies $X = YY^{\mathsf{T}}$ is a minimizer of (3.5).

5. The main algorithm. The main algorithm is summarized in Algorithm 5.1.

Algorithm 5.1 ManiSDP

```
Input: A, b, B, d, C, \sigma_0 > 0, \sigma_{\min} > 0, p_0 \in \mathbb{N}, \gamma > 1, \tau > 0, \delta_{ne} \in \mathbb{N}, (\epsilon_k)_{k \in \mathbb{N}} \subseteq \mathbb{R}^+
       with \lim_{k\to\infty} \epsilon_k = 0, \ \theta \in (0,1)
Output: (X, y, S)
  1: k \leftarrow 0, y^0 \leftarrow 0, p \leftarrow p_0, U \leftarrow 0_{n \times p};
  2: while stopping criteria do not fulfill do
           Solve the ALM subproblem (3.13) with U as a descent direction to obtain Y^{k+1}
           such that \|\operatorname{grad} \Psi_k(Y^{k+1})\|_F \leq \epsilon_k;

X^{k+1} \leftarrow Y^{k+1}(Y^{k+1})^{\mathsf{T}}, y^{k+1} \leftarrow y^k - \sigma_k(\mathcal{A}(X^{k+1}) - b);
           z_i \leftarrow \frac{\operatorname{Tr}(B_i X^{k+1} \nabla \Phi_k(X^{k+1}))}{\operatorname{Tr}(B_i^2 X^{k+1})}, i = 1, \dots, l;
  5:
           S^{k+1} \leftarrow C + \overset{\frown}{\sigma_k} \mathcal{A}^* (\overset{\frown}{\mathcal{A}}(X^{k+1}) - b - y^k / \sigma_k) - \mathcal{B}^*(z);
  6:
           Compute a descent direction U from S^{k+1} according to Theorem 4.4;
  7:
           Update p as described in Section 5.1;
  8:
           if \|A(X^{k+1}) - b\|/(1 + \|b\|) > \tau \|\operatorname{grad} \Psi_k(Y^{k+1})\|_F then
  9:
10:
               \sigma_{k+1} \leftarrow \gamma \sigma_k;
 11:
               \sigma_{k+1} \leftarrow \max \{ \sigma_k / \gamma, \sigma_{\min} \};
12:
           end if
13:
           k \leftarrow k+1;
14:
 15: end while
 16: return (X^k, y^k, S^k);
```

In Algorithm 5.1, the initial value p_0 can be typically set to 1 or 2, though for large-scale SDPs, a larger p_0 might be more advantageous; the value of γ is set to 2; the optimal choices for the values of the other parameters are problem-dependent. We discuss two important implementation details of Algorithm 5.1 in the subsequent subsections.

5.1. Dynamically adjusting the factorization size p. In implementation of Algorithm 5.1, the value of the factorization size p is dynamically adjusted. In particular, we increase p to escape from saddle points as discussed in Section 5.1; we decrease p to reduce the size of the ALM subproblem (3.13) so that the computational cost is as low as possible. Suppose that $\{s_i\}_i$ are the singular values of Y sorted from large to small. Then the rank of Y is estimated by

(5.1)
$$r = \arg\max_{i} \{i \mid s_i > \theta s_1\},$$

provided some threshold $\theta \in (0,1)$. Once the estimated rank r of Y is determined, we are able to construct a rank-r approximation of Y as follows. Suppose that Y has the singular value decomposition

$$Y = WDV^{\mathsf{T}},$$

where the diagonal of D is sorted from large to small. We then take

$$Y' = W_r D_r$$
.

as a rank-r approximation of Y, where W_r is the submatrix of W consisting of the first r columns and D_r is the upper-left $r \times r$ submatrix of D. Let

$$\delta = \min\{n_{\rm ne}, \delta_{\rm ne}\},\$$

where $n_{\rm ne}$ is the number of negative eigenvalues of S and $\delta_{\rm ne} \in \mathbb{N}$ is a tunable parameter. We then update the factorization size p by letting $p = r + \delta$ and accordingly let $Y = [Y', 0_{n \times \delta}]$. To obtain a descent direction, let $U = [0_{n \times r}, v_1, \cdots, v_{\delta}]$ where v_1, \ldots, v_{δ} are the eigenvectors corresponding to the δ smallest eigenvalues of S.

5.2. Self-adaptively updating the penalty parameter σ . Here we describe the strategy of self-adaptively updating the penalty parameter σ . Unlike usual ALMs using a monotonically nondecreasing sequence of penalty parameters, we allow to increase or decrease the penalty parameter self-adaptively. More concretely, we propose the following updating rules:

(5.2)
$$\sigma_{k+1} = \begin{cases} \gamma \sigma_k, & \text{if } \|\mathcal{A}(X^{k+1}) - b\|/(1 + \|b\|) > \tau \|\text{grad } \Psi_k(Y^{k+1})\|_F, \\ \max \{\sigma_k/\gamma, \sigma_{\min}\}, & \text{otherwise,} \end{cases}$$

where $\gamma > 1$, $\tau > 0$, $\sigma_{\min} > 0$ are constants. The intuition behind (5.2) is the following: the inequality $\|\mathcal{A}(X^{k+1}) - b\|/(1 + \|b\|) > \tau\|\operatorname{grad}\Psi_k(Y^{k+1})\|_F$ indicates that the progress of feasibility is not satisfactory and hence we increase the penalty parameter by setting $\sigma_{k+1} = \gamma \sigma_k$; otherwise, we decrease the penalty parameter by setting $\sigma_{k+1} = \max \{\sigma_k/\gamma, \sigma_{\min}\}$ so that the penalty parameter will not become too large or too small. We emphasis that a large penalty parameter may make the ALM subproblem (3.13) difficult to solve. Numerical experiments demonstrate that (5.2) works well in practice and significantly improves the performance of Algorithm 5.1.

5.3. Global convergence. We now give the global convergence result of Algorithm 5.1.

THEOREM 5.1. Assume that the sequence $(X^k)_{k\in\mathbb{N}}$ generated by Algorithm 5.1 satisfies $\lim_{k\to\infty}X^k=X^\star$. Assume that the sequences $(y^k)_{k\in\mathbb{N}}$, $(\Psi_k(Y^{k+1}))_{k\in\mathbb{N}}$ are bounded. Then X^\star is an optimal solution of (SDP-M). Moreover, let $\lim_{k\to\infty}y^k=y^\star$ and $\lim_{k\to\infty}S^k=S^\star$. Then $(X^\star,y^\star,S^\star)$ is a KKT point of (SDP-M), i.e., it satisfies (2.1) with z being given by

$$z_i \coloneqq \frac{\operatorname{Tr}(B_i X^*(C - \mathcal{A}^*(y^*)))}{\operatorname{Tr}(B_i^2 X^*)} \text{ for } i = 1, \dots, l.$$

Proof. We first show that for any threshold $\epsilon > 0$, there exists $K \in \mathbb{N}$ such that $\|\mathcal{A}(X^k) - b\| < \epsilon$ for all k > K. Suppose on the contrary that such K does not exist, i.e., $\|\mathcal{A}(X^k) - b\| \ge \epsilon$ for all sufficiently large k. Then Line 10 of Algorithm 5.1 is executed for infinitely many times and so $\lim_{k \to \infty} \sigma_k = +\infty$. Then note that

$$\Psi_k(Y^{k+1}) = \langle C, X^{k+1} \rangle + \frac{\sigma_k}{2} \left\| \mathcal{A}(X^{k+1}) - b - \frac{y^k}{\sigma_k} \right\|^2 - \frac{(y^k)^2}{2\sigma_k}$$

tends to infinity since $(y^k)_{k\in\mathbb{N}}$ is bounded, which contradicts to the assumption that $(\Psi_k(Y^{k+1}))_{k\in\mathbb{N}}$ is bounded. Therefore, $\|\mathcal{A}(X^*) - b\| = 0$ and X^* is a feasible solution of (SDP-M).

If X^* is not an optimal solution of (SDP-M), then by Theorem 3.10, we have $S^* \not\succeq 0$. So there exists a second-order descent direction for (3.13) starting from $Y^* = \lim_{k \to \infty} Y^k$ due to Theorem 4.4, which contradicts to the fact that X^* is the limit point of $(X^k)_{k \in \mathbb{N}}$.

The last assertion of the theorem can be easily verified.

6. Numerical experiments. In this section, we conduct comprehensive numerical experiments to benchmark our solver ManiSDP which implements Algorithm 5.1 in MATLAB. In particular, Manopt 7.1 [6] is employed by ManiSDP to solve the Riemannian manifold optimization problem (3.13). ManiSDP is freely available at https://github.com/wangjie212/ManiSDP-matlab.

Hardware. All numerical experiments were performed on a desktop computer with Intel(R) Core(TM) i9-10900 CPU@2.80GHz and 64G RAM.

Baseline Solvers. We compare the performance of ManiSDP with that of four advanced SDP solvers: MOSEK 10.0, SDPLR 1.03, SDPNAL+, STRIDE. We explain why to choose these four baseline solvers: MOSEK is chosen as it is a representative interior-point solver; SDPLR is chosen as it is a representative solver that also exploits the low-rank property via the Burer-Monteiro factorization; SDPNAL+ is chosen as it is a representative solver that combines first-order with second-order methods and is designed to solve large-scale SDPs; STRIDE is chosen as it specializes to solve large-scale SDP relaxations arising from polynomial optimization problems and exploits the rank-one property. In the following, running time of solvers is measured in seconds; "-" indicates that the solver encounters an out of memory error; "*" indicates that running time exceeds 10,000s; "**" indicates that the solver returns certain numerical error.

Stopping Criteria. To measure the feasibility and optimality of an approximate solution $(X, y, S) \in \mathbb{S}_n^+ \times \mathbb{R}^m \times \mathbb{S}_n^+$, we define the following residues:

(6.1)
$$\eta_p = \frac{\|\mathcal{A}(X) - b\|}{1 + \|b\|}, \ \eta_d = \frac{|\lambda_{\min}(S)|}{1 + |\lambda_{\max}(S)|}, \ \eta_g = \frac{|\langle C, X \rangle - b^{\mathsf{T}}y|}{1 + |\langle C, X \rangle| + |b^{\mathsf{T}}y|}.$$

Given a tolerance tol > 0, the SDP solver terminates when $\eta_{\text{max}} := \max \{\eta_p, \eta_d, \eta_g\} \le$ tol, and we set tol = 1e-8 for all our experiments.

Benchmark Problems. To benchmark the solvers, we solve six classes of SDPs arising from different situations (the Max-Cut problem, the matrix completion problem, binary quadratic programs, minimizing quartic polynomials on the unit sphere, the robust rotation search problem, nearest structured rank deficient matrices), with a focus on second-order SDP relaxations for polynomial optimization problems as they are highly degenerate and are challenging for most SDP solvers.

6.1. The Max-Cut problem. The Max-Cut problem is one of the basic combinatorial optimization problems, which is known to be NP-complete. Suppose that G(V, E) is an undirected graph with nodes $V = \{1, ..., N\}$ and with edge weights $w_{ij} = w_{ji}$ for $\{i, j\} \in E$. Then the Max-Cut problem for G, aiming to find the maximum cut, can be formulated as the following binary quadratic program:

(Max-Cut)
$$\begin{cases} \max & \frac{1}{2} \sum_{\{i,j\} \in E} w_{ij} (1 - x_i x_j) \\ \text{s.t.} & 1 - x_i^2 = 0, \quad i = 1, \dots, N. \end{cases}$$

To provide an upper bound on the maximum cut, we can consider the following SDP relaxation for (Max-Cut):

(6.2)
$$\begin{cases} \max & \frac{1}{4}\langle L, X \rangle \\ \text{s.t.} & X_{ii} = 1, \quad i = 1, \dots, N, \\ & X \succeq 0, \end{cases}$$

where $L \in \mathbb{S}_N$ is the Laplacian matrix of G, defined by

$$L_{ij} = \begin{cases} -w_{ij}, & \text{if } \{i, j\} \in E, \\ \sum_{k} w_{ik}, & \text{if } i = j, \\ 0, & \text{otherwise.} \end{cases}$$

Note that (6.2) fits in (SDP-M) with m = 0.

We select test graphs from the webpage https://web.stanford.edu/~yyye/yyye/Gset/ with N varying from 800 to 20000. For each instance, we solve (6.2) using the solvers MOSEK, SDPLR, SDPNAL+, and ManiSDP, respectively. The results are presented in Table 1. The following conclusions can be drawn from the table. (i) MOSEK can solve the instances with $N \leq 10000$ to high accuracy, but the running time significantly grows as N increases. When N = 20000, MOSEK runs out of space due to the large memory consumption of interior point methods. (ii) SDPNAL+ is very inefficient in solving this type of SDPs. For instance, when $N \geq 3000$, SDPNAL+ needs over 10000s to output the final result. (iii) Both SDPLR and ManiSDP can solve all instances to high accuracy, while ManiSDP is even more accurate and is faster than SDPLR by a factor of $2 \sim 10$.

6.2. The matrix completion problem. The matrix completion problem seeks to recover a low-rank matrix $M \in \mathbb{R}^{p \times q}$ from a subset of entries $\{M_{ij}\}_{(i,j)\in\Omega}$. This can be formulized as the convex optimization problem:

(MC)
$$\begin{cases} \inf_{Z \in \mathbb{R}^{p \times q}} & \|Z\|_* \\ \text{s.t.} & Z_{ij} = M_{ij}, \quad \forall (i, j) \in \Omega, \end{cases}$$

 $\begin{array}{c} {\rm TABLE~1} \\ {\it Results~for~the~Max-Cut~problem.} \end{array}$

graph	N	MOSEK	10.0	SDPLR	1.03	SDPN	AL+	Mani	ManiSDP		
grapii	1 V	$\eta_{ m max}$	time	/IIIax /IIIax		$\eta_{ m max}$	time	$\eta_{ m max}$	time		
G1	800	2.2e-09	3.51	7.4e-08	5.17	2.1e-09	52.2	1.6e-11	0.54		
G2	800	2.7e-09	3.54	2.1e-07	3.29	3.5e-09	52.5	5.3e-14	0.79		
G3	800	5.0e-09	3.50	2.1e-07	4.15	4.5e-09	38.6	4.3e-13	0.81		
G4	800	2.6e-09	3.41	2.3e-07	3.14	2.4e-09	37.9	9.8e-13	0.58		
G22	2000	1.0e-09	49.1	8.2e-08	12.5	8.6e-09	818	8.0e-12	1.48		
G23	2000	1.7e-09	51.3	2.7e-07	22.4	2.3e-08	555	5.8e-12	2.09		
G24	2000	9.2e-10	49.4	1.4e-06	6.95	4.3e-09	721	3.7e-12	1.36		
G25	2000	1.3e-09	53.8	3.6e-07	10.4	2.8e-09	770	1.9e-12	1.43		
G32	2000	2.3e-09	45.6	1.2e-07	22.9	7.0e-07	6463	1.6e-09	4.34		
G43	1000	4.0e-09	6.31	7.3e-08	2.47	3.0e-09	59.1	2.7e-13	0.68		
G44	1000	5.0e-09	6.27	3.1e-07	2.79	1.8e-08	61.2	4.2e-13	0.62		
G45	1000	1.1e-09	6.34	2.0e-07	2.73	1.6e-08	61.6	1.5e-12	0.59		
G48	3000	2.2e-09	108	1.1e-08	3.99	*	*	1.8e-17	1.81		
G49	3000	3.0e-10	100	4.6e-08	4.21	*	*	1.1e-16	1.94		
G50	3000	3.9e-14	112	3.0e-08	6.03	*	*	1.2e-14	2.25		
G55	5000	3.7e-09	963	2.3e-07	34.5	*	*	6.6e-12	17.7		
G56	5000	1.5e-09	847	7.5e-08	23.1	*	*	2.6e-12	15.4		
G57	5000	7.0e-10	877	1.5e-07	120	*	*	5.0e-09	30.8		
G58	5000	7.7e-10	1081	1.7e-07	101	*	*	1.2e-10	29.7		
G59	5000	4.5e-09	931	8.3e-08	63.6	*	*	2.3e-13	30.3		
G60	7000	1.1e-09	2722	7.5e-08	67.6	*	*	4.5e-12	35.9		
G61	7000	4.6e-10	2735	1.0e-07	114	*	*	4.4e-10	47.0		
G62	7000	2.2e-09	2484	3.1e-08	333	*	*	5.1e-09	124		
G63	7000	2.6e-09	2978	3.2e-07	224	*	*	9.6e-09	49.7		
G64	7000	1.2e-09	2886	5.8e-08	236	*	*	6.2e-09	51.9		
G65	8000	1.5e-09	3794	4.6e-08	307	*	*	1.2e-09	127		
G66	9000	2.6e-09	5464	5.7e-08	386	*	*	4.3e-09	169		
G67	10000	4.6e-09	7363	4.1e-08	610	*	*	7.0e-09	138		
G70	10000	4.1e-09	9451	3.1e-07	202	*	*	3.0e-12	73.3		
G72	10000	9.7e-11	7728	1.5e-07	614	*	*	8.3e-09	132		
G77	14000	*	*	6.9e-08	1177	*	*	1.9e-09	452		
G81	20000	-	-	5.2e-08	3520	*	*	8.3e-09	1934		

where $||Z||_* := \text{Tr}(Z^{\mathsf{T}}Z)^{\frac{1}{2}}$ is the nuclear norm of Z. Note that (MC) can be equivalently cast as an SDP of size $(n,m) = (p+q,|\Omega|)$:

(6.3)
$$\begin{cases} \inf_{X \in \mathbb{S}_n} & \operatorname{Tr}(X) \\ \text{s.t.} & \begin{bmatrix} 0_{p \times p} & E_{ij}^{\mathsf{T}} \\ E_{ij} & 0_{q \times q} \end{bmatrix} X = 2M_{ij}, \quad \forall (i, j) \in \Omega, \\ X = \begin{bmatrix} U & Z^{\mathsf{T}} \\ Z & V \end{bmatrix} \succeq 0, \end{cases}$$

where E_{ij} is a $p \times q$ matrix with 1 at its (i,j)-position and 0 otherwise. A famous result by Candes and Recht [10], later improved by Candes and Tao [11] states that, when M is low-rank and incoherent, and the number of samples satisfies $|\Omega| \geq Cn(\log n)^2$ with some constant C, then M can be exactly recovered by solving (6.3). In this subsection, we consider random instances of the matrix completion problem (MC). To this end, we select $\Omega \subseteq [p] \times [q]$ uniformly at random from all subsets with cardinality m, and set $M = M_1 M_2^{\mathsf{T}}$, where the entries of $M_1 \in \mathbb{R}^{p \times k}$ and $M_2 \in \mathbb{R}^{q \times k}$ are selected i.i.d. from the standard normal distribution. Here, we set k = 10, m = 400n and take p = q = 1000, 1500, 2000, 2500, 3000, 4000, 5000, 6000 respectively to generate test instances.

For each instance, we solve (6.3) using the solvers MOSEK, SDPLR, SDPNAL+, and ManiSDP, respectively. The results are presented in Table 2, from which we make the following observations. (i) MOSEK cannot solve any instance due to lack of enough memory. (ii) ManiSDP is not only the most efficient but also the most accurate among the remaining three solvers. Particularly, ManiSDP is faster than SDPLR by a factor of $1.5 \sim 2$, and is faster than SDPNAL+ by a order of magnitude.

6.3. Binary quadratic programs. Let us consider the binary quadratic program given by

(BQP)
$$\begin{cases} \inf_{\mathbf{x} \in \mathbb{R}^q} & \mathbf{x}^{\mathsf{T}} Q \mathbf{x} + \mathbf{c}^{\mathsf{T}} \mathbf{x} \\ \text{s.t.} & x_i^2 = 1, \quad i = 1, \dots, q, \end{cases}$$

where $Q \in \mathbb{S}_q$ and $\mathbf{c} \in \mathbb{R}^q$. (BQP) includes the Max-Cut problem (Max-Cut) as well as many other combinatorial optimization problems as special cases. On the other hand, (BQP) belongs to the more general class of polynomial optimization problems whose objective functions and constraints are given by polynomials. For a polynomial optimization problem, there is a systematic way to construct a hierarchy of increasingly tighter SDP relaxations, known as the moment-SOS hierarchy or the Lasserre hierarchy [22]. The moment SDP relaxation arising from the Lasserre hierarchy typically admits low-rank optimal solutions. Interestingly, for the binary quadratic program (BQP), the second-order moment relaxation is empirically tight on randomly generated instances as observed in [21, 39]. In the following we outline the ingredients of the second-order moment relaxation for (BQP). Let

$$v(\mathbf{x}) \coloneqq [1, x_1, \dots, x_q, x_1 x_2, x_1 x_3, \dots, x_{q-1} x_q]^\mathsf{T}$$

be the vector of monomials in \mathbf{x} up to degree two (excluding $x_i^2, i = 1, ..., q$) and $M := v(\mathbf{x})v(\mathbf{x})^{\mathsf{T}}$ be the corresponding moment matrix. Then the objective function

¹Under mild conditions, the optima of the hierarchy converge to the optimum of the polynomial optimization problem.

Table 2
Results for the matrix completion problem.

m	trial	m	MOSEK	10.0	SDPLR	1.03	SDPN	AL+	ManiSDP	
11	unan	111	0,536 0,590 0,590 0,0328 0,080 18,563 18,488 11,980 11,445 11,660 07,303 06,628 06,039 00,179 00,585 00,182 09,5929 - 96,602 96,604 993,420 - 94,532	time	$\eta_{ m max}$	time	$\eta_{ m max}$	time	$\eta_{ m max}$	time
	#1	550,536	-	-	1.7e-06	15.1	1.1e-08	69.9	2.1e-10	11.1
2000	#2	$550,\!565$	-	-	1.3e-06	14.7	4.5e-09	131	1.5e-10	10.7
	#3	$550,\!590$	-	-	8.6e-07	15.3	4.1e-09	143	7.4e-11	11.7
	#1	930,328	-	-	1.8e-06	51.4	3.1e-08	238	7.2e-10	28.0
3000	#2	929,882	-	-	6.7e-07	49.4	3.2e-08	217	3.6e-10	28.6
	#3	930,080	-	-	3.6e-06	45.4	3.1e-08	216	2.9e-10	28.3
,	#1	1,318,563	-	-	1.0e-06	88.7	4.8e-08	532	2.7e-10	57.0
4000	#2	1,318,488	-	-	1.6e-06	99.1	1.9e-09	548	1.4e-10	63.6
	#3	1,318,885	-	-	2.2e-06	96.8	4.7e-08	519	3.2e-10	61.7
	#1	1,711,980	-	-	1.2e-06	157	1.4e-09	1143	1.2e-10	104
5000	#2	1,711,445	-	-	1.0e-06	166	1.6e-09	1084	5.4e-11	103
	#3	1,711,660	-	-	1.1e-06	177	1.3e-09	1111	1.9e-10	104
,	#1	2,107,303	-	-	2.2e-07	272	2.1e-09	1883	2.5e-10	150
6000	#2	2,106,628	-	-	1.1e-06	260	2.2e-09	2001	2.8e-11	147
	#3	2,106,039	-	-	1.3e-06	271	2.5e-09	1979	1.3e-10	154
,	#1	2,900,179	-	-	2.1e-06	498	1.5e-08	3417	9.2e-11	249
8000	#2	2,900,585	-	-	3.4e-06	449	3.0e-08	4374	3.3e-10	259
	#3	2,900,182	-	-	3.2e-06	490	3.5e-08	4307	1.8e-10	248
	#1	3,695,929	-	-	1.1e-06	800	1.4e-09	8370	1.9e-10	446
10000	#2	3,696,602	-	-	2.1e-06	789	8.6e-09	8849	1.3e-10	456
	#3	3,696,604	-	-	1.1e-06	798	7.2e-09	8502	4.4e-10	472
	#1	4,493,420	-	-	7.8e-07	1310	*	*	2.6e-10	661
12000	#2	4,494,532	-	-	7.1e-07	1291	*	*	2.0e-10	696
	#3	4,493,391	-	-	3.5e-07	1330	*	*	2.2e-11	716

of (BQP) can be linearly expressed in terms of the entries of M. There are linear relationships among the entries of M consisting of $M_{ij} = M_{kr}$ whenever $M_{ij} - M_{kr}$ is reduced to 0 in the Gröbner basis $\{x_i^2 - 1\}_{i=1}^q$. Let $\mathcal{A}(X) = b$ collect all independent linear constraints obtained from these linear relationships when relaxing M to an unknown PSD matrix X. Moreover, because of the constraints $x_i^2 = 1, i = 1, \ldots, q$, the diagonal entries of M are all ones and so we let $\mathcal{B}(X) = d$ impose the unit-diagonal constraint on X. Consequently, we obtain the second-order moment relaxation for (BQP).

For each $q \in \{10, 20, 30, 40, 50, 60\}$, we generate three random instances of (BQP) by taking $Q \in \mathbb{S}_q$ with $Q_{ij} \sim \mathcal{N}(0,1)$ and $\mathbf{c} \in \mathbb{R}^q$ with $c_i \sim \mathcal{N}(0,1)$. For each instance, we solve the second-order moment relaxation using the solvers MOSEK, SDPLR, SDPNAL+, STRIDE and ManiSDP, respectively. The sizes of SDPs are recorded in Table 3 and the computational results are presented in Table 4. The following conclusions can be drawn from Table 4. (i) MOSEK can solve small-scale instances $(q \leq 20)$ to high accuracy, but the running time significantly grows as q increases (< 1s for q = 10 while ~ 50 s for q = 20). When $q \geq 30$, MOSEK runs out of space due to large memory consumption. (ii) SDPLR can solve small/medium-scale instances $(q \leq 30)$ to medium

Table 3
The sizes of SDPs for binary quadratic programs.

q	10	20	30	40	50	60
\overline{n}	56	211	466	821	1276	1831
\overline{m}	1,256	16,361	77,316	236,121	564,776	1,155,281

accuracy, but the running time significantly grows as q increases. When $q \geq 40$, SDPLR needs over 10000s to output the final result. (iii) SDPNAL+ can solve large-scale instances to medium/high accuracy, but the running time is pretty significant for large cases. (iv) Both STRIDE and ManiSDP can solve large-scale instances to high accuracy while ManiSDP is faster than STRIDE by a factor of $2 \sim 35$.

 $\begin{array}{c} {\rm TABLE} \ 4 \\ {\it Results \ for \ binary \ quadratic \ programs.} \end{array}$

a	trial	MOSEK	10.0	SDPLR	1.03	SDPN	AL+	STRI	DE	Manis	SDP		
q	unai	$\eta_{ m max}$	time										
	#1	2.6e-12	0.71	1.5e-06	0.52	1.9e-09	0.65	4.7e-13	0.79	3.9e-15	0.14		
10	#2	5.7e-14	0.84	6.0e-07	0.53	3.9e-09	1.37	3.4e-10	0.65	3.3e-15	0.18		
9 10 20 30 40 50	#3	8.0e-11	0.67	1.0e-06	1.27	1.5e-08	1.91	6.7e-13	0.68	4.2e-15	0.29		
	#1	9.8e-10	49.0	3.9e-07	30.8	3.0e-09	28.8	7.4e-13	6.12	1.5e-14	0.53		
20	#2	9.0e-10	50.3	2.3e-08	113	1.7e-08	29.0	6.4e-13	6.98	1.3e-14	0.61		
	#3	2.1e-12	47.9	6.6e-08	119	2.3e-07	12.5	2.9e-09	5.86	1.2e-14	0.72		
	#1	-	-	2.1e-06	8384	1.7e-04	187	1.2e-12	65.4	2.8e-14	3.93		
30	#2	-	-	2.7e-07	2796	6.4e-09	95.5	3.1e-09	36.2	3.2e-14	2.96		
	#3	-	-	1.6e-06	5698	7.8e-08	156	1.0e-12	60.3	2.8e-14	4.01		
	#1	-	-	*	*	2.1e-08	813	4.4e-13	249	4.6e-14	10.5		
40	#2	-	-	*	*	1.3e-06	1514	8.5e-09	294	4.7e-14	8.50		
	#3	-	-	*	*	1.3e-07	857	1.6e-12	321	4.4e-14	10.0		
	#1	-	-	*	*	1.6e-07	3058	7.8e-09	826	6.4e-14	31.1		
50	#2	-	-	*	*	4.8e-08	6347	1.8e-12	1020	8.9e-14	42.8		
	#3	-	-	*	*	7.0e-09	4800	8.2e-13	702	7.6e-14	61.4		
	#1	-	-	*	*	*	*	1.3e-12	2118	9.4e-14	94.3		
60	#2	-	-	*	*	*	*	*	*	9.5e-14	566		
	#3	-	-	*	*	*	*	3.3e-12	2704	8.7e-14	150		

To test the limit of ManiSDP, we run ManiSDP to solve the second-order moment relaxation of (BQP) with larger q. As shown in Table 5, ManiSDP can scale up to q = 120 for which the SDP has matrix dimension n = 7261 and contains m = 17,869,161 affine constraints, far beyond the reach of other SDP solvers!

6.4. Minimizing quartic polynomials on the unit sphere. Let us consider the problem of minimizing a quartic polynomial on the unit sphere:

(QS)
$$\begin{cases} \inf_{\mathbf{x} \in \mathbb{R}^q} \mathbf{c}^{\mathsf{T}} \cdot [\mathbf{x}]_4 \\ \text{s.t.} \quad \sum_{i=1}^q x_i^2 = 1, \end{cases}$$

where $[\mathbf{x}]_4$ is the vector of monomials in \mathbf{x} up to degree four and $\mathbf{c} \in \mathbb{R}^{|[\mathbf{x}]_4|}$. As for (BQP), the second-order moment relaxation is empirically tight on randomly gener-

q	70	80	90	100	110	120
n	2486	3241	4096	5051	6106	7261
m	2,119,636	3,589,841	5,717,896	8,675,801	12,655,556	17,869,161
$\eta_{ m max}$	1.4e-13	1.7e-13	2.2e-13	2.5e-13	3.0e-13	3.5e-13

 ${\it TABLE~5} \\ {\it Results~for~large~binary~quadratic~programs~via~ManiSDP}. \\$

ated instances of (QS) [39]. Let

$$v(\mathbf{x}) \coloneqq [1, x_1, \dots, x_q, x_1^2, x_1 x_2, x_1 x_3, \dots, x_{q-1} x_q, x_q^2]^\mathsf{T}$$

be the vector of monomials in \mathbf{x} up to degree two and $M := v(\mathbf{x})v(\mathbf{x})^{\mathsf{T}}$ be the corresponding moment matrix. Then the objective function of (QS) can be linearly expressed in terms of the entries of M. There are linear relationships among the entries of M consisting of all $M_{ij} = M_{kr}$. In addition, for each monomial $w \in v(\mathbf{x})$, the constraint $\sum_{i=1}^q x_i^2 = 1$ gives $w(\sum_{i=1}^q x_i^2 - 1) = 0$ which can be also linearly expressed in terms of the entries of M. Let A(X) = b collect all independent linear constraints obtained from these linear relationships when relaxing M to an unknown PSD matrix X. We therefore obtain the second-order moment relaxation for (QS).

For each $q \in \{10, 20, 30, 40, 50, 60\}$, we generate three random instances of (QS) by taking $\mathbf{c} \in \mathbb{R}^{|[\mathbf{x}]_4|}$ with $c_i \sim \mathcal{N}(0, 1)$. For each instance, we solve the second-order moment relaxation using the solvers MOSEK, SDPLR, SDPNAL+, STRIDE and ManiSDP, respectively. The results are presented in Table 6, from which we can draw the following conclusions. (i) MOSEK can solve small-scale instances ($q \leq 20$) to high accuracy, but the running time significantly grows as q increases (< 1s for q = 10 while ~ 50 s for q = 20). When $q \geq 30$, MOSEK runs out of space due to large memory consumption. (ii) SDPLR can solve all instances to medium accuracy, but the running time significantly grows as q increases. (iii) ManiSDP is the most efficient solver among the remaining three solvers. (iv) SDPNAL+ attains only medium accuracy for large-scale instances whereas STRIDE and ManiSDP can always attain high accuracy.

6.5. The robust rotation search problem. The robust rotation search problem (also known as the Wahba problem with outliers) is to search for the best 3D rotation to align two sets of 3D points while explicitly tolerating outliers, which can be formulated as the nonlinear optimization problem:

(6.4)
$$\min_{\|q\|=1} \sum_{i=1}^{N} \min \left\{ \frac{\|\tilde{z}_i - q \circ \tilde{w}_i \circ q^{-1}\|^2}{\beta_i^2}, 1 \right\},$$

where q is the unit quaternion parametrization of a 3D rotation, $(z_i \in \mathbb{R}^3, w_i \in \mathbb{R}^3)_{i=1}^N$ are given N pairs of 3D points, $\tilde{z} := [z^\intercal, 0]^\intercal \in \mathbb{R}^4$, $\tilde{w} := [w^\intercal, 0]^\intercal \in \mathbb{R}^4$, $q^{-1} := [-q_1, -q_2, -q_3, q_4]^\intercal$ is the inverse quaternion, "o" denotes the quaternion product, $\beta_i > 0$ is a given threshold that determines the maximum inlier residual. Problem (6.4) is a fundamental problem in aerospace, robotics and computer vision [37, 40]. By introducing N binary variables $\{\theta_i\}_{i=1}^N$, Problem (6.4) can be equivalently refor-

		MOSEK	10.0	SDPLR	1.03	SDPNA	SDPNAL+		DE	ManiSDP	
q	trial	$\eta_{ m max}$	time	$\eta_{ m max}$	time	$\eta_{ m max}$	time	$\eta_{ m max}$	time	$\eta_{ m max}$	time
	#1	6.6e-11	0.79	2.0e-07	0.07	2.5e-09	0.45	3.9e-12	0.35	4.5e-09	0.18
10	#2	8.8e-10	0.80	9.2e-07	0.56	1.3e-09	0.54	2.6e-12	0.52	7.0e-10	0.30
	#3	5.9e-10	0.79	9.0e-07	0.04	2.1e-09	0.54	2.3e-11	0.38	1.7e-10	0.18
	#1	7.4e-09	42.5	3.5e-06	1.35	1.2e-09	3.27	2.5e-12	2.74	3.7e-10	0.95
20	#2	4.0e-10	49.3	3.7e-07	2.34	5.6e-09	3.36	4.7e-11	2.91	5.4e-10	1.13
	#3	1.0e-08	42.9	5.2e-08	1.10	9.9e-09	3.20	8.7e-13	3.01	1.1e-09	0.82
	#1	-	-	1.5e-06	20.8	1.2e-09	20.9	3.1e-11	18.6	1.7e-10	6.46
30	#2	-	-	7.4e-07	38.4	1.5e-10	21.4	3.2e-13	19.4	8.4e-09	5.63
30	#3	-	-	1.8e-07	19.6	1.1e-09	19.0	1.1e-12	22.3	3.7e-10	5.92
	#1	-	-	3.5e-07	689	1.0e-07	45.0	2.8e-13	42.8	4.3e-09	28.7
40	#2	-	-	6.2e-07	272	1.4e-07	28.8	1.0e-12	39.1	4.2e-09	19.4
	#3	-	-	1.4e-07	261	3.9e-06	24.9	4.2e-11	39.3	8.9e-09	20.1
	#1	-	-	8.0e-07	1588	5.2e-07	68.9	2.3e-12	115	4.8e-09	61.4
50	#2	-	-	8.2e-08	1183	2.9e-06	69.0	8.4e-11	105	2.6e-09	49.9
	#3	-	-	3.9e-07	2350	1.1e-06	71.4	5.1e-11	124	4.7e-09	57.1
	#1	-	-	2.4e-07	4167	5.7e-07	177	2.6e-12	194	3.9e-09	109
60	#2	-	-	1.0e-08	7229	3.4e-07	237	3.6e-13	288	6.5e-10	116
	#3	-	-	3.6e-08	7752	4.7e-07	195	4.4e-13	209	2.1e-09	173

Table 6 Results for minimizing quartic polynomials on the unit sphere.

mulated as a polynomial optimization problem:

(RSS)
$$\min_{\substack{\|q\|=1,\\\theta_i \in \{-1,1\}, i=1,...,N}} \sum_{i=1}^{N} \frac{1+\theta_i}{2} \frac{\|\tilde{z}_i - q \circ \tilde{w}_i \circ q^{-1}\|^2}{\beta_i^2} + \frac{1-\theta_i}{2}.$$

Each θ_i acts as the selection variable for determining whether the i-th pair of 3D points (z_i, w_i) is an inlier or an outlier.

Yang and Carlone [37] proposed an SDP relaxation for (RSS) that was empirically shown to be tight. Let $\mathbf{x} = [q^{\intercal}, \theta_1, \dots, \theta_N]^{\intercal} \in \mathbb{R}^{N+4}$ be the decision variables of (RSS), and let

(6.5)
$$v(\mathbf{x}) = [q^{\mathsf{T}}, \theta_1 q^{\mathsf{T}}, \dots, \theta_N q^{\mathsf{T}}]^{\mathsf{T}} \in \mathbb{R}^{4N+4}$$

be the sparse set of monomials in \mathbf{x} of degree up to two. We build $M = v(\mathbf{x})v(\mathbf{x})^{\mathsf{T}}$ as the sparse moment matrix. Then the objective function of (RSS) can be linearly expressed in terms of the entries of M. There are linear relationships among the entries of M: (1) the diagonal 4×4 blocks of M are all identical $(\theta_i^2 q q^{\mathsf{T}} = q q^{\mathsf{T}})$; (2) the off-diagonal 4×4 blocks are symmetric $(\theta_i \theta_j q q^{\mathsf{T}} \in \mathbb{S}_4)$. Let $\mathcal{A}(X) = b$ collect all independent linear constraints obtained from these linear relationships when relaxing M to an unknown PSD matrix X. In addition, because of the unit quaternion constraint, M satisfies Tr(M) = N + 1 and so we let $\mathcal{B}(X) = d$ impose the trace constraint on X. Consequently, this leads to an SDP relaxation of size

(6.6)
$$n = 4N + 4, \quad m = 3N^2 + 13N + 1.$$

For each $N \in \{50, 100, 150, 200, 300, 500\}$, we generate three random instances of (RSS). For each instance, we solve the above SDP relaxation using the solvers MOSEK, SDPLR, SDPNAL+, STRIDE and ManiSDP, respectively. The results are presented in Table 7. The following conclusions can be drawn from the table. (i) MOSEK can solve small-scale instances ($N \leq 100$) to high accuracy, but the running time significantly grows as N increases ($N \leq 100$) to high accuracy, but the running time significantly grows as N increases ($N \leq 100$) while $N \leq 100$). When $N \geq 150$, MOSEK runs out of space due to large memory consumption. (ii) Both SDPLR and SDPNAL+ fail in solving these SDPs to even medium accuracy. (iii) Both STRIDE and ManiSDP can solve all instances to high accuracy while ManiSDP is faster than STRIDE by a factor of $N \sim 100$.

N	trial	MOSEK	10.0	SDPLR	1.03	SDPN	AL+	STRI	DE	Mani	SDP
11	ullai	$\eta_{ m max}$	time								
	#1	4.7e-10	16.4	9.8e-03	12.5	1.1e-02	106	2.8e-09	18.3	6.6e-09	3.02
50	#2	7.9e-10	19.3	3.1e-02	22.0	1.0e-02	96.3	7.3e-09	15.4	7.2e-10	2.93
	#3	1.1e-10	15.2	2.8e-03	19.4	1.1e-02	119	9.5e-09	15.4	5.5e-10	3.59
	#1	2.0e-11	622	3.6e-04	106	7.1e-02	642	3.1e-09	73.0	1.0e-09	22.9
100	#2	1.8e-10	653	8.1e-04	78.1	3.8e-02	631	1.6e-09	67.4	3.6e-10	20.3
	#3	7.3e-12	590	2.9e-03	67.2	7.8e-02	597	4.8e-09	69.0	5.4e-10	18.2
	#1	-	-	2.0e-03	291	8.0e-02	1691	4.3e-11	249	1.6e-09	33.5
150	#2	-	-	1.2e-03	233	6.4e-02	804	7.9e-09	171	4.7e-09	36.5
	#3	-	-	1.5e-01	416	1.2e-01	1491	9.9e-09	162	2.6e-09	33.8
	#1	-	-	3.1e-02	459	8.3e-02	2799	1.4e-09	254	6.3e-10	65.3
200	#2	-	-	1.6e-01	761	6.5e-02	1653	2.9e-09	306	9.2e-10	66.2
	#3	-	-	3.8e-03	894	6.3e-02	2171	3.2e-11	220	8.5e-10	67.9
	#1	-	-	1.1e-03	1264	5.2e-02	3528	4.1e-10	1176	1.1e-09	188
300	#2	-	-	7.3e-03	1787	4.9e-02	3421	8.0e-09	1458	3.6e-09	190
	#3	-	-	2.2e-03	1734	6.0e-02	4260	2.9e-09	868	1.2e-09	203
	#1	-	-	*	*	*	*	7.1e-09	5627	5.2e-10	601
500	#2	-	-	5.4e-02	9574	*	*	4.5e-10	4884	1.9e-09	801
	#3	_	_	*	*	*	*	3.4e-09	7878	5.0e-09	1055

Table 7
Results for the robust rotation search problem.

6.6. Nearest structured rank deficient matrices. Let us consider the problem of finding the nearest structured rank deficient matrix:

(6.7)
$$\min_{u \in \mathbb{R}^N} \left\{ \|u - \theta\|^2 \middle| L_0 + \sum_{i=1}^N u_i L_i \text{ is rank deficient} \right\},$$

where $L_i \in \mathbb{R}^{p \times q} (p \leq q), i = 0, ..., N$ and $\theta \in \mathbb{R}^N$ are given. Applications of Problem (6.7) (also known as the structured total least squares problem) could be found in [24]. We can reformulate (6.7) as the following polynomial optimization problem:

(NSRD)
$$\min_{z \in \mathbb{R}^p, u \in \mathbb{R}^N} \left\{ \|u - \theta\|^2 \middle| z^{\mathsf{T}} \left(L_0 + \sum_{i=1}^N u_i L_i \right) = 0, \|z\| = 1 \right\}.$$

Note that the unit vector z in (NSRD) serves as a witness of rank deficiency. (NSRD) is non-convex and Cifuentes proposed an SDP relaxation for (NSRD) [13] which is guaranteed to be tight under a low-noise assumption [14]. Let $\mathbf{x} = [z^{\mathsf{T}}, u^{\mathsf{T}}]^{\mathsf{T}} \in \mathbb{R}^{p+N}$

be the vector of variables involved in (NSRD), and let

(6.8)
$$v(\mathbf{x}) = [z^{\mathsf{T}}, u_1 z^{\mathsf{T}}, \dots, u_N z^{\mathsf{T}}]^{\mathsf{T}} \in \mathbb{R}^{p(N+1)}$$

be the sparse set of monomials in \mathbf{x} of degree up to two. We build $M = v(\mathbf{x})v(\mathbf{x})^{\mathsf{T}}$ as the sparse moment matrix. Then the objective function of (NSRD) can be linearly expressed in terms of the entries of M. There are linear relationships among the entries of M: (1) all off-diagonal $p \times p$ blocks $u_i u_j z z^{\mathsf{T}}$ are symmetric; (2) each of the first q equality constraint in (NSRD), say g = 0, gives rise to wg = 0 for each monomial $w \in v(\mathbf{x})$; (3) the unit norm of z implies that the trace of the leading $p \times p$ block of M is equal to 1. Let $\mathcal{A}(X) = b$ collect all independent linear constraints obtained from these linear relationships when relaxing M to an unknown PSD matrix X. Consequently, we obtain an SDP relaxation of size

$$n = p(N+1), \quad m = 1 + pq(N+1) + \frac{p(p-1)N(N+1)}{4}.$$

For each $p \in \{10, 15, 20, 25, 30, 40\}$, we generate three random instances of (6.7) with p=q and N=2p-1. For each instance, we solve the above SDP relaxation using the solvers MOSEK, SDPLR, SDPNAL+, STRIDE and ManiSDP, respectively. The results are presented in Table 8 from which we can make the following conclusions. (i) MOSEK can solve small-scale instances ($p \le 15$) to high accuracy, but the running time significantly grows as p increases ($p \le 15$) to high accuracy, but the running time significantly grows as p increases ($p \le 15$) to high accuracy, but the running time significantly grows as p increases ($p \le 15$) to high accuracy, but the running time significantly grows as p increases ($p \le 15$) to high accuracy, but the running time significantly grows as p increases ($p \le 15$) to high accuracy, and becomes unreliable when $p \ge 20$, MOSEK runs out of space due to large memory consumption. (ii) SDPLR can only solve small-scale instances to medium accuracy, and becomes unreliable when $p \ge 20$ for returning numerical errors. (iii) SDPNAL+ is much slower than STRIDE and ManiSDP, and can only obtain low/medium accuracy solutions. (iv) Both STRIDE and ManiSDP can solve the instances to high accuracy (occasionally STRIDE returns low/medium accuracy solutions) while ManiSDP is faster than STRIDE by a factor of $p \ge 10$.

7. Conclusions. We have presented a manifold optimization approach to solve linear SDPs with low-rank solutions by integrating the ALM and the Burer-Monteiro factorization. Global convergence is guaranteed despite of the non-convexity brought by the Burer-Monteiro factorization. Diverse numerical experiments demonstrate the superior performance of this approach. It has been shown that our solver ManiSDP is able to solve linear SDPs with millions of equality constraints to high accuracy in reasonable time.

We emphasis that ManiSDP is still at an early stage of development and the strength of the approach has not been fully revealed yet. Among others, we list several directions in enhancing the approach: (1) warm-starting; (2) using Riemannian line search to determine the step size for escaping from saddle points; (3) preconditioning for the Riemannian Hessian; (4) handling SDPs with inequality constraints. Moreover, as SDPs may contain multiple PSD blocks (e.g., SDP relaxations for sparse polynomial optimization problems [31, 32, 33]), it is also worth extending ManiSDP to handle multi-block SDPs. We believe that all of these efforts will eventually lead to a more powerful SDP solver, which makes large-scale low-rank SDPs even more tractable and hence allows to tackle hard application problems in real world.

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 ${\it TABLE~8} \\ {\it Results~for~nearest~structured~rank~deficient~matrices}. \\$

n	trial	MOSEK	10.0	SDPLR	1.03	SDPNA	AL+	STRI	DE	Mani	SDP
P	unai	$\eta_{ m max}$	time								
	#1	3.0e-11	22.9	8.4e-07	6.49	7.2e-08	64.1	3.5e-12	8.97	6.8e-10	1.28
10	#2	4.2e-11	20.1	6.4 e-05	3.04	1.8e-06	32.5	3.4e-12	4.74	4.4e-10	1.29
	#3	4.2e-09	15.3	6.1e-06	3.87	2.6e-05	15.5	1.3e-10	6.14	4.7e-09	0.90
	#1	4.9e-11	1623	1.5e-05	236	4.1e-06	233	4.4e-11	41.4	7.1e-09	12.7
15	#2	3.5e-09	1436	5.0e-05	369	2.9e-03	256	1.5e-10	33.0	7.2e-09	14.5
	#3	4.6e-10	1558	1.1e-05	32.5	1.8e-06	151	6.2e-11	35.5	6.5e-10	5.97
	#1	-	-	**	**	3.8e-03	894	3.0e-10	174	9.7e-09	55.9
20	#2	-	-	**	**	1.6e-02	1336	3.1e-11	125	7.9e-09	37.5
	#3	-	-	8.6e-06	1055	4.4e-03	1474	2.2e-10	149	7.5e-09	40.1
	#1	-	-	**	**	6.1e-03	8457	3.3e-06	4398	7.4e-09	781
25	#2	-	-	**	**	4.4e-07	3907	5.8e-10	429	7.9e-09	50.8
15	#3	-	-	**	**	1.3e-01	5153	2.6e-10	445	6.3e-09	75.5
	#1	-	-	*	*	*	*	4.2e-10	1812	4.9e-09	697
30	#2	-	-	*	*	*	*	4.2e-01	2484	7.2e-09	263
	#3	-	-	*	*	3.8e-07	9616	3.0e-11	1042	3.9e-09	108
_	#1	-	-	*	*	*	*	*	*	4.4e-09	1984
40	#2	-	-	*	*	*	*	*	*	4.0e-09	2493
	#3	-	-	*	*	*	*	*	*	3.3e-09	1279

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