Complete Facial Reduction in One Step for Spectrahedra

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Abstract

A spectrahedron is the feasible set of a semidefinite program, **SDP**, i.e., the intersection of an affine set with the positive semidefinite cone. While strict feasibility is a generic property for random problems, there are many classes of problems where strict feasibility fails and this means that strong duality can fail as well. If the minimal face containing the spectrahedron is known, the **SDP** can easily be transformed into an equivalent problem where strict feasibility holds and thus strong duality follows as well. The minimal face is fully characterized by the range or nullspace of any of the matrices in its relative interior. Obtaining such a matrix may require many facial reduction steps and is currently not known to be a tractable problem for spectrahedra with singularity degree greater than one. We propose a single parametric optimization problem with a resulting type of central path and prove that the optimal solution is unique and in the relative interior of the spectrahedron. Numerical tests illustrate the efficacy of our approach and its usefulness in regularizing **SDPs**.

Keywords: Semidefinite programming, SDP, facial reduction, singularity degree, maximizing log det.

AMS subject classifications: 90C22, 90C25

1 Introduction

A spectrahedron is the intersection of an affine manifold with the positive semidefinite cone

$$\mathcal{F} = \mathcal{F}(\mathcal{A}, b) := \{ X \in \mathbb{S}^n_+ : \mathcal{A}(X) = b \}, \tag{1.1}$$

where \mathbb{S}^n denotes the space of $n \times n$ symmetric matrices, $\mathbb{S}^n_+ \subset \mathbb{S}^n$ denotes the cone of positive semidefinite matrices, $\mathcal{A}: \mathbb{S}^n \to \mathbb{R}^m$ is a linear map, and $b \in \mathbb{R}^m$. Our motivation for studying spectrahedra arises from semidefinite programs, SDPs, where a linear objective is minimized over a spectrahedron. In contrast to e.g., linear programs, strong duality is not an inherent property of SDPs, but requires a constraint qualification (CQ) such as the Slater CQ, i.e., strict feasibility. For an SDP not satisfying the Slater CQ, the central path of the standard interior point algorithms is undefined and there is no guarantee of strong duality or convergence. Although instances where the Slater CQ fails are pathological, see e.g. [10] and [23], they occur in many applications and this phenomenon has lead to the development of a number of regularization methods, [7, 19, 20, 27, 28].

In this paper we focus on the *facial reduction* method, [3–5], where the optimization problem is restricted to the minimal face of \mathbb{S}^n_+ containing \mathcal{F} , denoted face(\mathcal{F}). Other regularization methods for **SDP** are given in [19, 20, 27, 28, 35]. See, for instance, the recent survey [9] for

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theory and applications. Further analyses and algorithms for facial reduction are discussed in [6, 22, 24–26, 34, 35].

The main structure of these approaches is that at each iteration a subproblem is solved to obtain an exposing vector for a face (not necessarily minimal) containing \mathcal{F} and the **SDP** is then reduced to this smaller face. The process is repeated until the **SDP** is reduced to face(\mathcal{F}). We highlight three challenges with this approach: (1) each subproblem is itself an **SDP** and thereby computationally intensive, (2) as many as n-1 subproblems may have to be solved, see [32], and (3) at each iteration a decision must be made regarding the rank of the exposing vector. In [6] it is shown that the subproblem satisfies the Slater CQ and in [25] a relaxation of the subproblem is given to ease the computational cost. A theoretical lower bound on the number of facial reduction steps, called the singularity degree, is introduced in [31] for algorithms relying on the theorem of the alternative (see Theorem 2.4). In [24] an alternative approach is given, using the self-dual embedding of [7] instead of the theorem of the alternative. This approach may not require as many steps as the singularity degree.

Our main contribution in this paper is a *primal* approach to facial reduction, which does not rely on exposing vectors, but instead obtains a matrix in the relative interior of \mathcal{F} , denoted relint(\mathcal{F}). Since the minimal face is characterized by the range of any such matrix, we obtain the facially reduced problem in just one step. As a result, we eliminate costly subproblems and require only one decision regarding rank.

While our motivation arises from **SDPs**, the problem of characterizing the relative interior of a spectrahedron is independent of this setting. The problem is formally stated below.

Problem 1.1. Given a spectrahedron $\mathcal{F}(A, b) \subseteq \mathbb{S}^n$, find $\bar{X} \in \text{relint}(\mathcal{F})$.

This paper is organized as follows. In Section 2 we introduce notation and discuss relevant material on **SDP** strong duality and facial reduction. We develop the theory for our approach in Section 3, prove convergence to the relative interior, and prove convergence to the analytic center under a sufficient condition. In Section 4, we propose an implementation of our approach and we present numerical results in Section 5. We also present a method for generating instances of **SDP** with varied singularity degree in Section 5.

2 Notation and Background

Throughout this paper the ambient space is the Euclidean space of $n \times n$ real symmetric matrices, \mathbb{S}^n , with the standard trace inner product $\langle X, Y \rangle := \operatorname{trace}(XY) = \sum_{i=1}^n \sum_{j=1}^n X_{ij} Y_{ij}$, and the induced Frobenius norm $||X||_F := \sqrt{\langle X, X \rangle}$. In the subsequent paragraphs, we highlight some well known results on the cone of positive semidefinite matrices and its faces, as well other useful results from convex analysis. For proofs and further reading we suggest [29, 32, 37].

The dimension of \mathbb{S}^n is the triangular number n(n+1)/2 =: t(n). We define svec: $\mathbb{S}^n \to \mathbb{R}^{t(n)}$ such that it maps the upper triangular elements of $X \in \mathbb{S}^n$ to a vector in $\mathbb{R}^{t(n)}$ where the off-diagonal elements are multiplied by $\sqrt{2}$. Then svec is an isometry and an isomorphism with $\mathrm{sMat} := \mathrm{svec}^{-1}$. Moreover, for $X, Y \in \mathbb{S}^n$, we have $\langle X, Y \rangle = \mathrm{svec}(X)^T \, \mathrm{svec}(Y)$. The eigenvalues of any $X \in \mathbb{S}^n$ are real and ordered so as to satisfy, $\lambda_1(X) \ge \cdots \ge \lambda_n(X)$, and $\lambda(X) \in \mathbb{R}^n$ is the vector consisting of all the eigenvalues. In terms of this notation, the operator 2-norm for matrices is defined as $\|X\|_2 := \max_i |\lambda_i(X)|$. When the argument to $\|\cdot\|_2$ is a vector, this denotes the usual Euclidean norm. The Frobenius norm may also be expressed in terms of eigenvalues: $\|X\|_F = \|\lambda(X)\|_2$. The set of positive semidefinite (PSD) matrices, \mathbb{S}^n_+ , is a closed convex cone in \mathbb{S}^n , whose interior consists of the positive definite (PD) matrices, \mathbb{S}^n_{++} . The cone \mathbb{S}^n_+ induces the $L\"{o}wner\ partial\ order\ on\ \mathbb{S}^n$. That is, for $X,Y\in\mathbb{S}^n$ we write $X\succeq Y$ when $X-Y\in\mathbb{S}^n_+$ and

similarly $X \succ Y$ when $X - Y \in \mathbb{S}_{++}^n$. Note that the following equivalence holds:

$$\langle X, Y \rangle = 0 \iff XY = 0, \ \forall X, Y \in \mathbb{S}^n_+.$$
 (2.1)

Definition 2.1 (face). A closed convex cone $f \subseteq \mathbb{S}^n_+$ is a face of \mathbb{S}^n_+ if

$$X, Y \in \mathbb{S}^n_+, X + Y \in f \implies X, Y \in f.$$

A nonempty face f is said to be *proper* if $f \neq \mathbb{S}^n_+$ and $f \neq 0$. Given a convex set $C \subseteq \mathbb{S}^n_+$, the *minimal face* of \mathbb{S}^n_+ containing f, with respect to set inclusion, is denoted face(C). A face f is said to be *exposed* if there exists $W \in \mathbb{S}^n_+ \setminus \{0\}$ such that

$$f = \{ X \in \mathbb{S}_+^n : \langle W, X \rangle = 0 \}.$$

Every face of \mathbb{S}^n_+ is exposed and the vector W is referred to as an *exposing vector*. The faces of \mathbb{S}^n_+ may be characterized in terms of the range of any of its maximal rank elements. Moreover, each face is isomorphic to a smaller dimensional positive semidefinite cone, as is seen in the subsequent theorem.

Theorem 2.2 ([9]). Let f be a face of \mathbb{S}^n_+ and $X \in f$ a maximal rank element with rank r and orthogonal spectral decomposition

$$X = \begin{bmatrix} V & U \end{bmatrix} \begin{bmatrix} D & 0 \\ 0 & 0 \end{bmatrix} \begin{bmatrix} V & U \end{bmatrix}^T \in \mathbb{S}^n_+, \quad D \in \mathbb{S}^r_{++}.$$

Then $f = V\mathbb{S}_+^r V^T$ and $\operatorname{relint}(f) = V\mathbb{S}_{++}^r V^T$. Moreover, $W \in \mathbb{S}_+^n$ is an exposing vector for f if and only if $W \in U\mathbb{S}_{++}^{n-r} U^T$.

We refer to $U\mathbb{S}^{n-r}_+U^T$, from the above theorem, as the *conjugate face*, denoted f^c . For any convex set C, an explicit form for face(C) and face(C)^c may be obtained from the orthogonal spectral decomposition of any of its maximal rank elements as in Theorem 2.2.

For a linear map $\mathcal{A}: \mathbb{S}^n \to \mathbb{R}^m$, there exist $S_1, \ldots, S_m \in \mathbb{S}^n$ such that

$$(\mathcal{A}(X))_i = \langle X, S_i \rangle, \quad \forall i \in \{1, \dots, m\}.$$

The adjoint of \mathcal{A} is the unique linear map $\mathcal{A}^*: \mathbb{R}^m \to \mathbb{S}^n$ satisfying

$$\langle \mathcal{A}(X), y \rangle = \langle X, \mathcal{A}^*(y) \rangle, \quad \forall X \in \mathbb{S}^n, \ y \in \mathbb{R}^m,$$

and has the explicit form $\mathcal{A}^*(y) = \sum_{i=1}^m y_i S_i$, i.e., range $(\mathcal{A}^*) = \text{span}\{S_1, \dots, S_m\}$. We define $A_i \in \mathbb{S}^n$ to form a basis for the nullspace, null $(\mathcal{A}) = \text{span}\{A_1, \dots, A_q\}$.

For a non-empty convex set $C \subseteq \mathbb{S}^n$ the recession cone, denoted C^{∞} , captures the directions in which C is unbounded. That is

$$C^{\infty} := \{ Y \in \mathbb{S}^n : X + \lambda Y \in C, \ \forall \lambda \ge 0, \ X \in C \}.$$

$$(2.2)$$

Note that the recession directions are the same at all points $X \in C$. For a non-empty set $S \subseteq \mathbb{S}^n$, the dual cone (also referred to as the positive polar) is defined as

$$S^{+} := \{ Y \in \mathbb{S}^{n} : \langle X, Y \rangle > 0, \ \forall X \in S \}.$$
 (2.3)

For cones K_1 and K_2 , with $cl(\cdot)$ denoting set closure, a useful result regarding dual cones is:

$$(K_1 \cap K_2)^+ = \operatorname{cl}(K_1^+ + K_2^+), \tag{2.4}$$

2.1 Strong Duality in Semidefinite Programming and Facial Reduction Consider the standard primal form SDP

$$\mathbf{SDP} \qquad p^* := \min\{\langle C, X \rangle : \mathcal{A}(X) = b, X \succeq 0\}, \tag{2.5}$$

with Lagrangian dual

D-SDP
$$d^* := \min\{b^T y : \mathcal{A}^*(y) \leq C\}. \tag{2.6}$$

Let \mathcal{F} denote the spectrahadron defined by the feasible set of \mathbf{SDP} . One of the challenges in semidefinite programming is that strong duality is not an inherent property, but depends on a constraint qualification, such as the Slater \mathbf{CQ} .

Theorem 2.3 (strong duality, [37]). If the primal optimal value p^* is finite and $\mathcal{F} \cap \mathbb{S}^n_{++} \neq \emptyset$, then the primal-dual pair SDP and D-SDP have a zero duality gap, $p^* = d^*$, and d^* is attained.

Since the Lagrangian dual of the dual is the primal, this result can similarly be applied to the dual problem, i.e., if the primal-dual pair both satisfy the Slater CQ, then there is a zero duality gap and both optimal values are attained.

Not only can strong duality fail with the absence of the Slater CQ, but the standard central path of an interior point algorithm is undefined. The facial reduction regularization approach of [3–5] restricts **SDP** to the minimal face of \mathbb{S}^n_+ containing \mathcal{F} :

SDP-R
$$\min\{\langle C, X \rangle : \mathcal{A}(X) = b, X \in \text{face}(\mathcal{F})\}.$$
 (2.7)

Since the dimension of \mathcal{F} and face(\mathcal{F}) is the same, the Slater CQ holds for the facially reduced problem. Moreover, face(\mathcal{F}) is isomorphic to a smaller dimensional positive semidefinite cone, thus **SDP-R** is itself a semidefinite program. The restriction to face(\mathcal{F}) may be obtained as in the results of Theorem 2.2. The dual of **SDP-R** restricts the slack variable to the dual cone

$$Z = C - \mathcal{A}^*(y) \in \text{face}(\mathcal{F})^+.$$

Note that $\mathcal{F}^+ = \mathrm{face}(\mathcal{F})^+$. If we have knowledge of $\mathrm{face}(\mathcal{F})$, i.e., we have the matrix V such that $\mathrm{face}(\mathcal{F}) = V\mathbb{S}_+^r V^T$, then we may replace X in \mathbf{SDP} by VRV^T with $R \succeq 0$. After rearranging, we obtain \mathbf{SDP} - \mathbf{R} . Alternatively, if our knowledge of the minimal face is in the form of an exposing vector, say W, then we may obtain V so that its columns form a basis for $\mathrm{null}(W)$. We see that the approach is straightforward when knowledge of $\mathrm{face}(\mathcal{F})$ is available. In instances where such knowledge is unavailable, the following theorem of the alternative from [5] guarantees the existence of exposing vectors that lie in $\mathrm{range}(\mathcal{A}^*)$.

Theorem 2.4 (of the alternative, [5]). Exactly one of the following systems is consistent:

1.
$$A(X) = b, X \succ 0,$$

2.
$$0 \neq A^*(y) \succeq 0, b^T y = 0.$$

The first alternative is just the Slater CQ, while if the second alternative holds, then $\mathcal{A}^*(y)$ is an exposing vector for a face containing \mathcal{F} . We may use a basis for null($\mathcal{A}^*(y)$) to obtain a smaller \mathbf{SDP} . If the Slater CQ holds for the new \mathbf{SDP} we have obtained $\mathbf{SDP-R}$, otherwise, we find an exposing vector and reduce the problem again. We outline the facial reduction procedure in Algorithm 2.1. At each iteration, the dimension of the problem is reduced by at least one, hence this approach is bound to obtain $\mathbf{SDP-R}$ in at most n-1 iterations, assuming that the initial problem is feasible. If at each iteration the exposing vector obtained is of maximal rank then the

Algorithm 2.1 Facial reduction procedure using the theorem of the alternative.

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Initialize S_i so that (\mathcal{A}(X))_i = \langle S_i, X \rangle for i \in \{1, \dots, m\} while Item 2. of Theorem 2.4 do obtain exposing vector W W = \begin{bmatrix} U & V \end{bmatrix} \begin{bmatrix} D & 0 \\ 0 & 0 \end{bmatrix} \begin{bmatrix} U & V \end{bmatrix}, \quad D \succ 0 S_i \leftarrow V^T S_i V, \quad i \in \{1, \dots, m\} end while
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number of iterations required to obtain **SDP-R** is referred to as the *singularity degree*, [31]. For a non-empty spectrahedron, \mathcal{F} , we denote the singularity degree as $sd = sd(\mathcal{F})$.

We remark that any algorithm pursuing the minimal face through exposing vectors of the form $\mathcal{A}^*(\cdot)$, must perform at least as many iterations as the singularity degree. The singularity degree could be as large as the trivial upper bound n-1 as is seen in the example of [32]. Thus facial reduction may be very expensive computationally. On the other hand, from Theorem 2.2 we see that face(\mathcal{F}) is fully characterized by the range of any of its relative interior matrices. That is, from any solution to Problem 1.1 we may obtain the regularized problem **SDP-R**.

3 A Parametric Optimization Approach

In this section we present a parametric optimization problem that solves Problem 1.1.

Assumption 3.1. We make the following assumptions: (1) A is surjective; (2) F is non-empty, bounded and contained in a proper face of \mathbb{S}^n_+ .

The assumption on \mathcal{A} is a standard regularity assumption and so is the non-emptiness assumption on \mathcal{F} . The necessity of \mathcal{F} to be bounded will become apparent throughout this section, however, our approach may be applied to unbounded spectrahedra as well. We discuss such extensions in Section 3.2. The assumption that \mathcal{F} is contained in a proper face of \mathbb{S}^n_+ restricts our discussion to those instances of **SDP** that are interesting with respect to facial reduction.

In the following lemma are stated two useful characterizations of bounded spectrahedra.

Lemma 3.2. The following holds:

$$\mathcal{F}$$
 is bounded \iff null $(\mathcal{A}) \cap \mathbb{S}^n_+ = \{0\} \iff \operatorname{range}(\mathcal{A}^*) \cap \mathbb{S}^n_{++} \neq \emptyset$.

Proof. For the first equivalence, \mathcal{F} is bounded if and only if $\mathcal{F}^{\infty} = \{0\}$ by Theorem 8.4 of [29]. It suffices, therefore, to show that $\mathcal{F}^{\infty} = \text{null}(\mathcal{A}) \cap \mathbb{S}^n_+$. It is easy to see that $(\mathbb{S}^n_+)^{\infty} = \mathbb{S}^n_+$ and that the recession cone of the affine manifold defined by \mathcal{A} and b is $\text{null}(\mathcal{A})$. By Corollary 8.3.3 of [29] the recession cone of the intersection of convex sets is the intersection of the respective recession cones, yielding the desired result.

Now let us consider the second equivalence. For the forward direction, observe that

$$\operatorname{null}(\mathcal{A}) \cap \mathbb{S}_{+}^{n} = \{0\} \iff \left(\operatorname{null}(\mathcal{A}) \cap \mathbb{S}_{+}^{n}\right)^{+} = \{0\}^{+},$$
$$\iff \operatorname{null}(\mathcal{A})^{\perp} + \mathbb{S}_{+}^{n} = \mathbb{S}^{n},$$
$$\iff \operatorname{range}(\mathcal{A}^{*}) + \mathbb{S}_{+}^{n} = \mathbb{S}^{n}.$$

The second inequality is due to (2.4) and one can verify that in this case $\operatorname{null}(\mathcal{A})^{\perp} \cap \mathbb{S}_{+}^{n}$ is closed. Thus there exists $X \in \operatorname{range}(\mathcal{A}^{*})$ and $Y \in \mathbb{S}_{+}^{n}$ such that X + Y = -I. Equivalently, $-X = I + Y \in \mathbb{S}_{++}^{n}$. For the converse, let $X \in \operatorname{range}(\mathcal{A}^{*}) \cap \mathbb{S}_{++}^{n}$ and suppose $0 \neq S \in \operatorname{null}(\mathcal{A}) \cap \mathbb{S}_{+}^{n}$. Then $\langle X, S \rangle = 0$ which implies, by (2.1), that XS = 0. But then $\operatorname{null}(X) \neq \{0\}$, a contradiction. \square

Let r denote the maximal rank of any matrix in relint(\mathcal{F}) and let the columns of $V \in \mathbb{R}^{n \times r}$ form a basis for its range. In seeking a relative interior point of \mathcal{F} we define a specific point from which we develop a parametric optimization problem.

Definition 3.3 (analytic center). The analytic center of \mathcal{F} is the unique matrix \hat{X} satisfying

$$\hat{X} = \arg\max\{\log\det(V^T X V) : X \in \mathcal{F}\}. \tag{3.1}$$

Under Assumption 3.1 the analytic center is well-defined and this follows from the proof of Theorem 3.4, below. It is easy to see that the analytic center is indeed in the relative interior of \mathcal{F} and therefore a solution to Probelm 1.1. However, the optimization problem from which it is derived is intractable due to the unknown matrix V. If V is simply removed from the optimization problem (replaced with the identity), then the problem is ill-posed since the objective does not take any finite values over the feasible set as it lies on the boundary of the **SDP** cone. To combat these issues, we propose replacing V with I and also perturbing \mathcal{F} so that it intersects \mathbb{S}^n_{++} . The perturbation we choose is that of replacing b with $b(\alpha) := b + \alpha \mathcal{A}(I)$, $\alpha > 0$, thereby defining a family of spectrahedra

$$\mathcal{F}(\alpha) := \{ X \in \mathbb{S}^n_+ : \mathcal{A}(X) = b(\alpha) \}.$$

It is easy to see that if $\mathcal{F} \neq \emptyset$ then $\mathcal{F}(\alpha)$ has postive definite elements for every $\alpha > 0$. Indeed $\mathcal{F} + \alpha I \subset \mathcal{F}(\alpha)$. Note that the affine manifold may be perturbed by any positive definite matrix and I is chosen for simplicity. We now consider the family of optimization problems for $\alpha > 0$:

$$\mathbf{P}(\alpha) \qquad \max\{\log \det(X) : X \in \mathcal{F}(\alpha)\}. \tag{3.2}$$

It is well known that the solution to this problem exists and is unique for each $\alpha > 0$. We include a proof in Theorem 3.4, below. Moreover, since $face(\mathcal{F}(\alpha)) = \mathbb{S}^n_+$ for each $\alpha > 0$, the solution to $\mathbf{P}(\alpha)$ is in relint($\mathcal{F}(\alpha)$) and is exactly the analytic center of $\mathcal{F}(\alpha)$. The intuition behind our approach is that as the perturbation gets smaller, i.e., $\alpha \searrow 0$, the solution to $\mathbf{P}(\alpha)$ approaches the relative interior of \mathcal{F} . This intuition is validated in Section 3.3. Specifically, we show that the solutions to $\mathbf{P}(\alpha)$ form a smooth path that converges to $\bar{X} \in \text{relint}(\mathcal{F})$. We also provide a sufficient condition for the limit point to be \hat{X} in Section 3.4.

We note that our approach of perturbing the spectrahedron in order to use the $\log \det(\cdot)$ function is not entirely new. In [12], for instance, the authors perturb a convex feasible set in order to approximate the rank function using $\log \det(\cdot)$. Unlike our approach, their perturbation is constant.

3.1 Optimality Conditions

We choose the strictly concave function $\log \det(\cdot)$ for its elegant optimality conditions, though the maximization is equivalent to maximizing only the determinant. We treat it as an *extended* valued concave function that takes the value $-\infty$ if X is singular. For this reason we refer to both functions $\det(\cdot)$ and $\log \det(\cdot)$ equivalently throughout our discussion.

Let us now consider the optimality conditions for the problem $\mathbf{P}(\alpha)$. Similar problems have been thoroughly studied throughout the literature in matrix completions and \mathbf{SDP} , e.g., [1, 14, 33, 37]. Nonetheless, we include a proof for completeness and to emphasize its simplicity.

Theorem 3.4 (optimality conditions). For every $\alpha > 0$ there exists a unique $X(\alpha) \in \mathcal{F}(\alpha) \cap \mathbb{S}^n_{++}$ such that

$$X(\alpha) = \arg\max\{\log\det(X) : X \in \mathcal{F}(\alpha)\}. \tag{3.3}$$

Moreover, $X(\alpha)$ satisfies (3.3) if, and only if, there exists a unique $y(\alpha) \in \mathbb{R}^m$ and a unique $Z(\alpha) \in \mathbb{S}^n_{++}$ such that

$$\begin{bmatrix} \mathcal{A}^*(y(\alpha)) - Z(\alpha) \\ \mathcal{A}(X(\alpha)) - b(\alpha) \\ Z(\alpha)X(\alpha) - I \end{bmatrix} = 0.$$
 (3.4)

Proof. By Assumption 3.1, $\mathcal{F} \neq \emptyset$ and bounded and it follows that $\mathcal{F}(\alpha) \cap \mathbb{S}^n_{++} \neq \emptyset$ and by Lemma 3.2 it is bounded. Moreover, $\log \det(\cdot)$ is a strictly concave function over $\mathcal{F}(\alpha) \cap \mathbb{S}^n_{++}$ (a so-called barrier function) and

$$\lim_{\det(X)\to 0} \log \det(X) = -\infty.$$

Thus, we conclude that the optimum $X(\alpha) \in \mathcal{F}(\alpha) \cap \mathbb{S}^n_{++}$ exists and is unique. The Lagrangian of problem (3.3) is

$$\mathcal{L}(X, y) = \log \det(X) - \langle y, \mathcal{A}(X) - b \rangle$$

= \log \det(X) - \langle \mathcal{A}^*(y), X \rangle + \langle y, b \rangle.

Since the constraints are linear, stationarity of the Lagrangian holds at $X(\alpha)$. Hence there exists $y(\alpha) \in \mathbb{R}^m$ such that $(X(\alpha))^{-1} = \mathcal{A}^*(y(\alpha)) =: Z(\alpha)$. Clearly $Z(\alpha)$ is unique, and since \mathcal{A} is surjective, we conclude in addition that $y(\alpha)$ is unique.

3.2 The Unbounded Case

Before we continue with the convergence results, we briefly address the case of unbounded spectrahedra. The restriction to bounded spectrahedra is necessary in order to have solutions to (3.3). There are certainly large families of **SDPs** where the assumption holds. Problems arising from liftings of combinatorial optimization problems often have the diagonal elements specified, and hence bound the corresponding spectrahedron. Matrix completion problems are another family where the diagonal is often specified. Nonetheless, many **SDPs** have unbounded feasible sets and we provide two methods for reducing such spectrahedra to bounded ones. First, we show that the boundedness of \mathcal{F} may be determined by solving a projection problem.

Proposition 3.5. Let \mathcal{F} be a spectrahedron defined by the affine manifold $\mathcal{A}(X) = b$ and let

$$P := \arg\min \{ ||X - I||_F : X \in \operatorname{range}(\mathcal{A}^*) \}.$$

Then \mathcal{F} is bounded if $P \succ 0$.

Proof. First we note that P is well defined and a singleton since it is the projection of I onto a closed convex set. Now $P \succ 0$ implies that $\operatorname{range}(\mathcal{A}^*) \cap \mathbb{S}^n_{++} \neq \emptyset$ and by Lemma 3.2 this is equivalent to \mathcal{F} bounded.

The proposition gives us a sufficient condition for \mathcal{F} to be bounded. Suppose this condition is not satisfied, but we have knowledge of some matrix $S \in \mathcal{F}$. Then for t > 0, consider the spectrahedron

$$\mathcal{F}' := \{ X \in \mathbb{S}^n : X \in \mathcal{F}, \operatorname{trace}(X) = \operatorname{trace}(S) + t \}.$$

Clearly \mathcal{F}' is bounded. Moreover, we see that $\mathcal{F}' \subset \mathcal{F}$ and contains maximal rank elements of \mathcal{F} , hence face(\mathcal{F}') = face(\mathcal{F}). It follows that relint(\mathcal{F}') \subset relint(\mathcal{F}) and we have reduced the problem to the bounded case.

Now suppose that the sufficient condition of the proposition does not hold and we do not have knowledge of a feasible element of F. In this case we detect recession directions, elements

of $\text{null}(\mathcal{A}) \cap \mathbb{S}^n_+$, and project to the orthogonal complement. Specifically, if \mathcal{F} is unbounded then $\mathcal{F}(\alpha)$ is unbounded and problem (3.2) is unbounded. Suppose, we have detected unboundedness, i.e., we have $X \in \mathcal{F}(\alpha) \cap \mathbb{S}^n_+$ with large norm. Then $X = S_0 + S$ with $S \in \text{null}(\mathcal{A}) \cap \mathbb{S}^n_+$ and $||S|| \gg ||S_0||$. We then restrict \mathcal{F} to the orthogonal complement of S, that is, we consider the new spectrahedron

$$\mathcal{F}' := \{ X \in \mathbb{S}^n : X \in \mathcal{F}, \ \langle S, X \rangle = 0 \}.$$

By repeated application, we eliminate a basis for the recession directions and obtain a bounded spectrahedron. From any of the relative interior points of this spectrahedron, we may obtain a relative interior point for \mathcal{F} by adding to it the recession directions obtained throughout the reduction process.

3.3 Convergence to the Relative Interior and Smoothness

By simple inspection it is easy to see that $(X(\alpha), y(\alpha), Z(\alpha))$, as in (3.4), does not converge as $\alpha \searrow 0$. Indeed, under Assumption 3.1,

$$\lim_{\alpha \searrow 0} \lambda_n(X(\alpha)) \to 0 \implies \lim_{\alpha \searrow 0} ||Z(\alpha)||_2 \to +\infty.$$

It is therefore necessary to scale $Z(\alpha)$ so that it remains bounded. Let us look at an example.

Example 3.6. Consider the matrix completion problem: find $X \succeq 0$ having the form

$$\begin{pmatrix} 1 & 1 & ? \\ 1 & 1 & 1 \\ ? & 1 & 1 \end{pmatrix}.$$

The set of solutions is indeed a spectrahedron with A and b given by

$$\mathcal{A}\left(\begin{bmatrix} x_{11} & x_{12} & x_{13} \\ x_{12} & x_{22} & x_{23} \\ x_{13} & x_{23} & x_{33} \end{bmatrix}\right) := \begin{pmatrix} x_{11} \\ x_{12} \\ x_{22} \\ x_{23} \\ x_{33} \end{pmatrix}, b := \begin{pmatrix} 1 \\ 1 \\ 1 \\ 1 \\ 1 \end{pmatrix}.$$

In this case, it is not difficult to obtain

$$X(\alpha) = \begin{pmatrix} 1 + \alpha & 1 & \frac{1}{1+\alpha} \\ 1 & 1 + \alpha & 1 \\ \frac{1}{1+\alpha} & 1 & 1+\alpha \end{pmatrix},$$

with inverse

$$X(\alpha)^{-1} = \frac{1}{\alpha(2+\alpha)} \begin{pmatrix} 1+\alpha & -1 & 0\\ -1 & \frac{\alpha^2+2\alpha+2}{1+\alpha} & -1\\ 0 & -1 & 1+\alpha \end{pmatrix}.$$

Clearly $\lim_{\alpha\searrow 0} ||X(\alpha)^{-1}||_2 \to +\infty$. However, when we consider $\alpha X(\alpha)^{-1}$, and take the limit as α goes to 0 we obtain the bounded limit

$$\bar{Z} = \begin{pmatrix} \frac{1}{2} & -\frac{1}{2} & 0\\ -\frac{1}{2} & 1 & -\frac{1}{2}\\ 0 & -\frac{1}{2} & \frac{1}{2} \end{pmatrix}.$$

Note that $\bar{X} = X(0)$ is the 3×3 matrix with all ones, $\operatorname{rank} \bar{X} + \operatorname{rank} \bar{Z} = 3$, and $\bar{X}\bar{Z} = 0$.

It turns out that multiplying $X(\alpha)^{-1}$ by α always bounds the sequence $(X(\alpha), y(\alpha), Z(\alpha))$. Therefore, we consider the scaled system

$$\begin{bmatrix} \mathcal{A}^*(y) - Z \\ \mathcal{A}(X) - b(\alpha) \\ ZX - \alpha I \end{bmatrix} = 0, \ X \succ 0, \ Z \succ 0, \ \alpha > 0, \tag{3.5}$$

that is obtained from (3.4) by multiplying the last equation by α . Abusing our previous notation, we let $(X(\alpha), y(\alpha), Z(\alpha))$ denote a solution to *this* system and we refer to the set of all such solutions as the *parametric path*. The parametric path has clear parallels to the *central path* of **SDP**, however, it differs in one main respect: it is not contained in the relative interior of \mathcal{F} . In the main theorems of this section we prove that the parametric path is smooth and converges as $\alpha \searrow 0$ with the primal limit point in relint(\mathcal{F}). We begin by showing that the primal component of the parametric path has cluster points.

Lemma 3.7. Let $\bar{\alpha} > 0$. For every sequence $\{\alpha_k\}_{k \in \mathbb{N}} \subset (0, \bar{\alpha}]$ such that $\alpha_k \searrow 0$, there exists a subsequence $\{\alpha_l\}_{l \in \mathbb{N}}$ such that $X(\alpha_l) \to \bar{X} \in \mathcal{F}$.

Proof. Let $\bar{\alpha}$ and $\{\alpha_k\}_{k\in\mathbb{N}}$ be as in the hypothesis. First we show that the sequence $X(\alpha_k)$ is bounded. For any $k\in\mathbb{N}$ we have

$$||X(\alpha_k)||_2 \le ||X(\alpha_k) + (\bar{\alpha} - \alpha_k)I||_2 \le \max_{X \in \mathcal{F}(\bar{\alpha})} ||X||_2 < +\infty.$$

The second inequality is due to $X(\alpha_k) + (\bar{\alpha} - \alpha_k)I \in \mathcal{F}(\bar{\alpha})$ and the third inequality holds since $\mathcal{F}(\bar{\alpha})$ is bounded. Thus there exists a convergent subsequence $\{\alpha_l\}_{l\in\mathbb{N}}$ with $X(\alpha_l) \to \bar{X}$, that clearly belongs to \mathcal{F} .

For the dual variables we need only prove that $Z(\alpha)$ converges (for a subsequence) since this implies that $y(\alpha)$ also converges, by the assumption that \mathcal{A} is surjective. As for $X(\alpha)$, we show that the tail of the parametric path corresponding to $Z(\alpha)$ is bounded. To this end, we first prove the following technical lemma. Recall that \hat{X} is the analytic center of Definition 3.3.

Lemma 3.8. Let $\bar{\alpha} > 0$. There exists M > 0 such that for all $\alpha \in (0, \bar{\alpha}]$,

$$0 < \langle X(\alpha)^{-1}, \hat{X} + \alpha I \rangle \le M.$$

Proof. Let $\bar{\alpha}$ be as in the hypothesis and let $\alpha \in (0, \bar{\alpha}]$. The first inequality is trivial since both of the matrices are positive definite. For the second inequality, we have,

$$\langle X(\bar{\alpha})^{-1} - X(\alpha)^{-1}, \hat{X} + \bar{\alpha}I - X(\alpha) \rangle = \langle \frac{1}{\bar{\alpha}} \mathcal{A}^*(y(\bar{\alpha})) - \frac{1}{\alpha} \mathcal{A}^*(y(\alpha)), \hat{X} + \bar{\alpha}I - X(\alpha) \rangle,$$

$$= \langle \frac{1}{\bar{\alpha}} y(\bar{\alpha}) - \frac{1}{\alpha} y(\alpha), \mathcal{A}(\hat{X} + \bar{\alpha}I) - \mathcal{A}(X(\alpha)) \rangle,$$

$$= \langle \frac{1}{\bar{\alpha}} y(\bar{\alpha}) - \frac{1}{\alpha} y(\alpha), (\bar{\alpha} - \alpha) \mathcal{A}(I) \rangle,$$

$$= \langle X(\bar{\alpha})^{-1} - X(\alpha)^{-1}, (\bar{\alpha} - \alpha)I \rangle,$$

$$= (\bar{\alpha} - \alpha) \operatorname{trace}(X(\bar{\alpha})^{-1}) - \langle X(\alpha)^{-1}, (\bar{\alpha} - \alpha)I \rangle.$$

$$(3.6)$$

On the other hand,

$$\langle X(\bar{\alpha})^{-1} - X(\alpha)^{-1}, \hat{X} + \bar{\alpha}I - X(\alpha) \rangle = n + \langle X(\bar{\alpha})^{-1}, \hat{X} \rangle + \bar{\alpha}\operatorname{trace}(X(\bar{\alpha})^{-1}) - \langle X(\bar{\alpha})^{-1}, X(\alpha) \rangle - \langle X(\alpha)^{-1}, \hat{X} + \bar{\alpha}I \rangle.$$
(3.7)

Combining (3.6) and (3.7) we get

$$(\bar{\alpha} - \alpha)\operatorname{trace}(X(\bar{\alpha})^{-1}) - \langle X(\alpha)^{-1}, (\bar{\alpha} - \alpha)I \rangle = n + \langle X(\bar{\alpha})^{-1}, \hat{X} \rangle + \bar{\alpha}\operatorname{trace}(X(\bar{\alpha})^{-1}) - \langle X(\bar{\alpha})^{-1}, X(\alpha) \rangle - \langle X(\alpha)^{-1}, \hat{X} + \bar{\alpha}I \rangle.$$

After rearranging, we obtain

$$\langle X(\alpha)^{-1}, \hat{X} + \alpha I \rangle = n + \langle X(\bar{\alpha})^{-1}, \hat{X} \rangle + \bar{\alpha} \operatorname{trace}(X(\bar{\alpha})^{-1}) - \langle X(\bar{\alpha})^{-1}, X(\alpha) \rangle - (\bar{\alpha} - \alpha) \operatorname{trace}(X(\bar{\alpha})^{-1}),$$

$$= n + \alpha \operatorname{trace}(X(\bar{\alpha})^{-1}) + \langle X(\bar{\alpha})^{-1}, \hat{X} \rangle - \langle X(\bar{\alpha})^{-1}, X(\alpha) \rangle.$$
(3.8)

The first and the third terms of the right hand side are positive constants. The second term is positive for every value of α and is bounded above by $\bar{\alpha} \operatorname{trace}(X(\bar{\alpha})^{-1})$ while the fourth term is bounded above by 0. Applying these bounds as well as the trivial lower bound on the left hand side, we get

$$0 < \langle X(\alpha)^{-1}, \hat{X} + \alpha I \rangle \le n + \bar{\alpha} \operatorname{trace}(X(\bar{\alpha})^{-1}) + \langle X(\bar{\alpha})^{-1}, \hat{X} \rangle =: M.$$
(3.9)

We need one more ingredient to prove that the parametric path corresponding to $Z(\alpha)$ is bounded. This involves bounding the trace inner product above and below by the maximal and minimal scalar products of the eigenvalues, respectively.

Lemma 3.9 (Ky-Fan [11], Hoffman-Wielandt [17]). If $A, B \in \mathbb{S}^n$, then

$$\sum_{i=1}^{n} \lambda_i(A)\lambda_{n+1-i}(B) \le \langle A, B \rangle \le \sum_{i=1}^{n} \lambda_i(A)\lambda_i(B).$$

We now have the necessary tools for proving boundedness and obtain the following convergence result.

Theorem 3.10. Let $\bar{\alpha} > 0$. For every sequence $\{\alpha_k\}_{k \in \mathbb{N}} \subset (0, \bar{\alpha}]$ such that $\alpha_k \searrow 0$, there exists a subsequence $\{\alpha_\ell\}_{\ell \in \mathbb{N}}$ such that

$$(X(\alpha_{\ell}), y(\alpha_{\ell}), Z(\alpha_{\ell})) \to (\bar{X}, \bar{y}, \bar{Z}) \in \{\mathbb{S}_{+}^{n} \times \mathbb{R}^{m} \times \mathbb{S}_{+}^{n}\}$$

with $\bar{X} \in \operatorname{relint}(\mathcal{F})$ and $\bar{Z} = \mathcal{A}^*(\bar{y})$.

Proof. Let $\bar{\alpha} > 0$ and $\{\alpha_k\}_{k \in \mathbb{N}}$ be as in the hypothesis. We may without loss of generality assume that $X(\alpha_k) \to \bar{X} \in \mathcal{F}$ due to Lemma 3.7. Let $k \in \mathbb{N}$. Combining the upper bound of Lemma 3.8 with the lower bound of Lemma 3.9 we have $\sum_{i=1}^n \lambda_i (X(\alpha_k)^{-1}) \lambda_{n+1-i} (\hat{X} + \alpha_k I) \leq M$. Since the left hand side is a sum of positive terms, the inequality applies to each term:

$$\lambda_i(X(\alpha_k)^{-1})\lambda_{n+1-i}(\hat{X} + \alpha_k I) \le M, \quad \forall i \in \{1, \dots, n\}.$$

Equivalently,

$$\lambda_i(X(\alpha_k)^{-1}) \le \frac{M}{\lambda_{n+1-i}(\hat{X}) + \alpha_k}, \quad \forall i \in \{1, \dots, n\}.$$
(3.10)

Now exactly r eigenvalues of \hat{X} are positive. Thus for $i \in \{n-r+1,\ldots,n\}$ we have

$$\lambda_i(X(\alpha_k)^{-1}) \le \frac{M}{\lambda_{n+1-i}(\hat{X}) + \alpha_k} \le \frac{M}{\lambda_{n+1-i}(\hat{X})},$$

and we conclude that the r smallest eigenvalues of $X(\alpha_k)^{-1}$ are bounded above. Consequently, there are at least r eigenvalues of $X(\alpha_k)$ that are bounded away from 0 and $\operatorname{rank}(\bar{X}) \geq r$. On the other hand $\bar{X} \in \mathcal{F}$ and $\operatorname{rank}(\bar{X}) \leq r$ and it follows that $\bar{X} \in \operatorname{relint}(\mathcal{F})$.

Now we show that $Z(\alpha_k)$ is a bounded sequence. Indeed, from (3.10) we have

$$||Z(\alpha_k)||_2 = \alpha_k \lambda_1(X(\alpha_k)^{-1}) \le \alpha_k \frac{M}{\lambda_n(\hat{X}) + \alpha_k} = \alpha_k \frac{M}{\alpha_k} = M.$$

The second to last equality follows from the assumption that $\hat{X} \in \mathbb{S}^n_+ \setminus \mathbb{S}^n_{++}$, i.e. $\lambda_n(\hat{X}) = 0$. Now there exists a subsequence $\{\alpha_\ell\}_{\ell \in \mathbb{N}}$ such that $Z(\alpha_\ell) \to \bar{Z}$, $X(\alpha_\ell) \to \bar{X}$. Moreover, for each ℓ , there exists a unique $y(\alpha_\ell) \in \mathbb{R}^m$ such that $Z(\alpha_\ell) = \mathcal{A}^*(y(\alpha_\ell))$ and since \mathcal{A} is surjective, there exists $\bar{y} \in \mathbb{R}^m$ such that $y(\alpha_\ell) \to \bar{y}$ and $\bar{Z} = \mathcal{A}^*(\bar{y})$. Lastly, the sequence $Z(\alpha_\ell)$ is contained in the closed cone \mathbb{S}^n_+ hence $\bar{Z} \in \mathbb{S}^n_+$, completing the proof.

We conclude this section by proving that the parametric path is smooth and has a limit point as $\alpha \searrow 0$. Our proof relies on the following lemma of Milnor and is motivated by an analogous proof for the central path of **SDP** in [15,16]. Recall that an *algebraic set* is the solution set of a system of finitely many polynomial equations.

Lemma 3.11 (Milnor [21]). Let $\mathcal{V} \subseteq \mathbb{R}^k$ be an algebraic set and $\mathcal{U} \subseteq \mathbb{R}^k$ be an open set defined by finitely many polynomial inequalities. Then if $0 \in \operatorname{cl}(\mathcal{U} \cap \mathcal{V})$ there exists $\varepsilon > 0$ and a real analytic curve $p : [0, \varepsilon) \to \mathbb{R}^k$ such that p(0) = 0 and $p(t) \in \mathcal{U} \cap \mathcal{V}$ whenever t > 0.

Theorem 3.12. There exists $(\bar{X}, \bar{y}, \bar{Z}) \in \mathbb{S}^n_+ \times \mathbb{R}^m \times \mathbb{S}^n_+$ with all the properties of Theorem 3.10 such that

$$\lim_{\alpha \searrow 0} (X(\alpha), y(\alpha), Z(\alpha)) = (\bar{X}, \bar{y}, \bar{Z}).$$

Proof. Let $(\bar{X}, \bar{y}, \bar{Z})$ be a cluster point of the parametric path as in Theorem 3.10. We define the set \mathcal{U} as

$$\mathcal{U} := \{ (X, y, Z, \alpha) \in \mathbb{S}^n \times \mathbb{R}^m \times \mathbb{S}^n \times \mathbb{R} : \bar{X} + X \succ 0, \ \bar{Z} + Z \succ 0, \ Z = \mathcal{A}^*(y), \ \alpha > 0 \}.$$

Note that each of the positive definite constraints is equivalent to n strict determinant (polynomial) inequalities. Therefore, \mathcal{U} satisfies the assumptions of Lemma 3.11. Next, let us define the set \mathcal{V} as,

$$\mathcal{V} := \left\{ (X, y, Z, \alpha) \in \mathbb{S}^n \times \mathbb{R}^m \times \mathbb{S}^n \times \mathbb{R} : \begin{bmatrix} \mathcal{A}^*(y) - Z \\ \mathcal{A}(X) + \alpha \mathcal{A}(I) \\ (\bar{Z} + Z)(\bar{X} + X) - \alpha I \end{bmatrix} = 0 \right\},$$

and note that \mathcal{V} is indeed a real algebraic set. Next we show that there is a one-to-one correspondance between $\mathcal{U} \cap \mathcal{V}$ and the parametric path without any of its cluster points. Consider $(\tilde{X}, \tilde{y}, \tilde{Z}, \tilde{\alpha}) \in \mathcal{U} \cap \mathcal{V}$ and let $(X(\tilde{\alpha}), y(\tilde{\alpha}), Z(\tilde{\alpha}))$ be a point on the parametric path. We show that

$$(\bar{X} + \tilde{X}, \bar{y} + \tilde{y}, \bar{Z} + \tilde{Z}) = (X(\tilde{\alpha}), y(\tilde{\alpha}), Z(\tilde{\alpha})). \tag{3.11}$$

First of all $\bar{X} + \tilde{X} \succ 0$ and $\bar{Z} + \tilde{Z} \succ 0$ by inclusion in \mathcal{U} . Secondly, $(\bar{X} + \tilde{X}, \bar{y} + \tilde{y}, \bar{Z} + \tilde{Z})$ solves the system (3.5) when $\alpha = \tilde{\alpha}$:

$$\begin{bmatrix} \mathcal{A}^*(\bar{y}+\tilde{y}) - (\bar{Z}+\tilde{Z}) \\ \mathcal{A}(\bar{X}+\tilde{X}) - b(\tilde{\alpha}) \\ (\bar{Z}+\tilde{Z})(\bar{X}+\tilde{X}) - \tilde{\alpha}I \end{bmatrix} = \begin{bmatrix} \mathcal{A}^*(\bar{y}) - \bar{Z} + (\mathcal{A}^*(\tilde{y}) - \tilde{Z}) \\ b + \tilde{\alpha}\mathcal{A}(I) - b(\tilde{\alpha}) \\ 0 \end{bmatrix} = \begin{bmatrix} 0 \\ 0 \\ 0 \end{bmatrix}.$$

Since (3.5) has a unique solution, (3.11) holds. Thus,

$$(\tilde{X}, \tilde{y}, \tilde{Z}) = (X(\alpha) - \bar{X}, y(\alpha) - \bar{y}, Z(\alpha) - \bar{Z}),$$

and it follows that $\mathcal{U} \cap \mathcal{V}$ is a translation of the parametric path (without its cluster points):

$$\mathcal{U} \cap \mathcal{V} = \{ (X, y, Z, \alpha) \in \mathbb{S}^n \times \mathbb{R}^m \times \mathbb{S}^n \times \mathbb{R} : (X, y, Z) = (X(\alpha) - \bar{X}, y(\alpha) - \bar{y}, Z(\alpha) - \bar{Z}), \ \alpha > 0 \}.$$
(3.12)

Next, we show that $0 \in \operatorname{cl}(\mathcal{U} \cap \mathcal{V})$. To see this, note that

$$(X(\alpha), y(\alpha), Z(\alpha)) \to (\bar{X}, \bar{y}, \bar{Z}),$$

as $\alpha \setminus 0$ along a subsequence. Therefore, along the same subsequence, we have

$$(X(\alpha) - \bar{X}, y(\alpha) - \bar{y}, Z(\alpha) - \bar{Z}, \alpha) \to 0.$$

Each of the elements of this subsequence belongs to $\mathcal{U} \cap \mathcal{V}$ by (3.12) and therefore $0 \in \operatorname{cl}(\mathcal{U} \cap \mathcal{V})$. We have shown that \mathcal{U} and \mathcal{V} satisfy all the assumptions of Lemma 3.11, hence there exists $\varepsilon > 0$ and an analytic curve $p : [0, \varepsilon) \to \mathbb{S}^n \times \mathbb{R}^m \times \mathbb{S}^n \times \mathbb{R}$ such that p(0) = 0 and $p(t) \in \mathcal{U} \cap \mathcal{V}$ for t > 0. Let

$$p(t) = (X_{(t)}, y_{(t)}, Z_{(t)}, \alpha_{(t)}),$$

and observe that by (3.12), we have

$$(X_{(t)}, y_{(t)}, Z_{(t)}, \alpha_{(t)}) = (X(\alpha_{(t)}) - \bar{X}, y(\alpha_{(t)}) - \bar{y}, Z(\alpha_{(t)}) - \bar{Z}). \tag{3.13}$$

Since p is a real analytic curve, the map $g:[0,\varepsilon)\to\mathbb{R}$ defined as $g(t)=\alpha_{(t)}$, is a differentiable function on the open interval $(0,\varepsilon)$ with

$$\lim_{t \searrow 0} g(t) = 0.$$

In particular, this implies that there is an interval $[0,\bar{\varepsilon}) \subseteq [0,\varepsilon)$ where g is monotone. It follows that on $[0,\bar{\varepsilon})$, g^{-1} is a well defined continuous function that converges to 0 from the right. Note that for any t > 0, (X(t), y(t), Z(t)) is on the parametric path. Therefore,

$$\lim_{t \searrow 0} X(t) = \lim_{t \searrow 0} X(g(g^{-1}(t))) = \lim_{t \searrow 0} X(\alpha_{(g^{-1}(t))}).$$

Substituting with (3.13), we have

$$\lim_{t \searrow 0} X(t) = \lim_{t \searrow 0} X_{(g^{-1}(t))} + \bar{X} = \bar{X}.$$

Similarly, y(t) and Z(t) converge to \bar{y} and \bar{Z} respectively. Thus every cluster point of the parametric path is identical to $(\bar{X}, \bar{y}, \bar{Z})$.

We have shown that the tail of the parametric path is smooth and it has a limit point. Smoothness of the entire path follows from the Berge Maximum Theorem, [2], or [30, Example 5.22].

3.4 Convergence to the Analytic Center

The results of the previous section establish that the parametric path converges to relint(\mathcal{F}) and therefore the primal part of the limit point has excatly r positive eigenvalues. If the smallest positive eigenvalue is very small it may be difficult to distinguish it from zero numerically. Therefore it is desirable for the limit point to be *substantially* in the relative interior, in the sense that its smallest positive eigenvalue is relatively large. The analytic center has this property and so a natural question is whether the limit point coincides with the analytic center. In the following modification of an example of [16], the parametric path converges to a point different from the analytic center.

Example 3.13. Consider the SDP feasibility problem where A is defined by the matrices

and $b := (1,0,0,0,0)^T$. One can verify that the feasible set consists of positive semidefinite matrices of the form

and the analytic center is the determinant maximizer over the positive definite blocks of this set and satisfies $x_{22} = 0.5$ and $x_{12} = 0$. However, the parametric path converges to a matrix with $x_{22} = 0.6$ and $x_{12} = 0$. To see this note that

$$\mathcal{A}(I) = \begin{pmatrix} 2 & 1 & 1 & 0 & 1 \end{pmatrix}^T, \quad b(\alpha) = \begin{pmatrix} 1 + 2\alpha & \alpha & \alpha & 0 & \alpha \end{pmatrix}^T.$$

By feasibility, $X(\alpha)$ has the form

$$\begin{bmatrix} 1 + 2\alpha - x_{22} & x_{12} & x_{13} & x_{14} \\ x_{12} & x_{22} & 0 & \frac{1}{2}(\alpha - x_{33}) \\ x_{13} & 0 & x_{33} & 0 \\ x_{14} & \frac{1}{2}(\alpha - x_{33}) & 0 & \alpha \end{bmatrix}.$$

Moreover, the optimality conditions of Theorem 3.4 indicate that $X(\alpha)^{-1} \in \text{range}(\mathcal{A}^*)$ and

$$X(\alpha) = \begin{bmatrix} 1 + 2\alpha - x_{22} & 0 & 0 & 0\\ 0 & x_{22} & 0 & \frac{1}{2}(\alpha - x_{33})\\ 0 & 0 & x_{33} & 0\\ 0 & \frac{1}{2}(\alpha - x_{33}) & 0 & \alpha \end{bmatrix}.$$

Of all the matrices with this form, $X(\alpha)$ is the one maximizing the determinant, that is

$$([X(\alpha)]_{22}, [X(\alpha)]_{33})^T = \underset{s.t.}{\arg\max} \quad x_{33}(1 + 2\alpha - x_{22})(\alpha x_{22} - \frac{1}{4}(\alpha - x_{33})^2), \\ s.t. \quad 0 < x_{22} < 1 + 2\alpha, \, x_{33} > 0, \, \alpha x_{22} > \frac{1}{4}(\alpha - x_{33})^2.$$

Due to the strict inequalities, the maximizer is a stationary point of the objective function. Computing the derivative with respect to x_{22} and x_{33} we obtain the equations

$$x_{33}(-(\alpha x_{22} - \frac{1}{4}(\alpha - x_{33})^2) + \alpha(1 + 2\alpha - x_{22}) = 0,$$

$$(1 + 2\alpha - x_{22})((\alpha x_{22} - \frac{1}{4}(\alpha - x_{33})^2) + \frac{1}{2}x_{33}(\alpha - x_{33})) = 0.$$

Since $x_{33} > 0$ and $(1 + 2\alpha - x_{22}) > 0$, we may divide them out. Then solving each equation for x_{22} we get

$$x_{22} = \frac{1}{8\alpha}(\alpha - x_{33})^2 + \alpha + \frac{1}{2} = \frac{1}{4\alpha}(\alpha - x_{33})^2 - \frac{1}{2\alpha}x_{33}(\alpha - x_{33}).$$
 (3.14)

Combining these two expressions for x_{22} we have

$$0 = \frac{1}{4\alpha}(\alpha - x_{33})^2 - \frac{1}{2\alpha}x_{33}(\alpha - x_{33}) - \frac{1}{8\alpha}(\alpha - x_{33})^2 - \alpha - \frac{1}{2},$$

$$= \frac{1}{8\alpha}(\alpha - x_{33})^2 - \frac{1}{2}x_{33} + \frac{1}{2\alpha}x_{33}^2 - \alpha - \frac{1}{2},$$

$$= \frac{1}{8\alpha}x_{33}^2 - \frac{1}{4}x_{33} + \frac{1}{8}\alpha - \frac{1}{2}x_{33} + \frac{1}{2\alpha}x_{33}^2 - \alpha - \frac{1}{2},$$

$$= \frac{5}{8\alpha}x_{33}^2 - \frac{3}{4}x_{33} + \frac{1}{8}\alpha - \alpha - \frac{1}{2}.$$

Now we solve for x_{33} ,

$$x_{33} = \frac{\frac{3}{4} \pm \sqrt{\frac{9}{16} - 4(\frac{5}{8\alpha})(\frac{1}{8}\alpha - \alpha - \frac{1}{2})}}{2\frac{5}{8\alpha}} = \frac{3\alpha}{5} \pm \frac{4\alpha}{5}\sqrt{\frac{11\alpha + 5}{4\alpha}} = \frac{1}{5}(3\alpha + 2\sqrt{\alpha}\sqrt{11\alpha + 5}).$$

Since x_{33} is fully determined by the stationarity constraints, we have $[X(\alpha)]_{33} = x_{33}$ and $[X(\alpha)]_{33} \to 0$ as $\alpha \searrow 0$. Substituting this expression for x_{33} into (3.14) we get

$$\begin{split} [X(\alpha)]_{22} &= \frac{1}{8\alpha}(\alpha - \frac{1}{5}(3\alpha + 2\sqrt{\alpha}\sqrt{11\alpha + 5}))^2 + \alpha + \frac{1}{2}, \\ &= \frac{1}{8\alpha}(\alpha^2 - 2\alpha\frac{1}{5}(3\alpha + 2\sqrt{\alpha}\sqrt{11\alpha + 5}) + \frac{1}{25}(9\alpha^2 + 6\alpha\sqrt{\alpha}\sqrt{11\alpha + 5} + 4\alpha(11\alpha + 5))) + \alpha + \frac{1}{2}, \\ &= \frac{1}{8}\alpha - \frac{1}{20}(3\alpha + 2\sqrt{\alpha}\sqrt{11\alpha + 5}) + \frac{1}{200}(9\alpha + 6\sqrt{\alpha}\sqrt{11\alpha + 5} + 4(11\alpha + 5)) + \alpha + \frac{1}{2}, \\ &= \frac{31}{25}\alpha - \frac{7}{100}\sqrt{\alpha}\sqrt{11\alpha + 5} + \frac{6}{10}. \end{split}$$

Now it is clear that $[X(\alpha)]_{22} \to 0.6$ as $\alpha \searrow 0$.

3.4.1 A Sufficient Condition for Convergence to the Analytic Center

Recall that face $(\mathcal{F}) = V \mathbb{S}_+^r V^T$. To simplify the discussion we may assume that $V = \begin{bmatrix} I \\ 0 \end{bmatrix}$, so that

$$face(\mathcal{F}) = \begin{bmatrix} \mathbb{S}_{+}^{r} & 0\\ 0 & 0 \end{bmatrix}. \tag{3.15}$$

This follows from the rich automorphism group of \mathbb{S}^n_+ , that is, for any full rank $W \in \mathbb{R}^{n \times n}$, we have $W\mathbb{S}^n_+W^T=\mathbb{S}^n_+$. Moreover, it is easy to see that there is a one-to-one correspondence between relative interior points under such transformations.

Let us now express \mathcal{F} in terms of null(\mathcal{A}), that is, if $A_0 \in \mathcal{F}$ and recall that $A_1, \ldots, A_q, q = t(n) - m$, form a basis for null(\mathcal{A}), then

$$\mathcal{F} = (A_0 + \operatorname{span}\{A_1, \dots, A_q\}) \cap \mathbb{S}_+^n.$$

Similarly,

$$\mathcal{F}(\alpha) = (\alpha I + A_0 + \operatorname{span}\{A_1, \dots, A_q\}) \cap \mathbb{S}_+^n.$$

Next, let us partition A_i according to the block structure of (3.15):

$$A_i = \begin{bmatrix} L_i & M_i \\ M_i^T & N_i \end{bmatrix}, \quad i \in \{0, \dots, q\}.$$

$$(3.16)$$

Since $A_0 \in \mathcal{F}$, from (3.15) we have $N_0 = 0$ and $M_0 = 0$. Much of the subsequent discussion focuses on the linear pencil $\sum_{i=1}^{q} x_i N_i$. Let \mathcal{N} be the linear mapping such that

$$\operatorname{null}(\mathcal{N}) = \left\{ \sum_{i=1}^{q} x_i N_i : x \in \mathbb{R}^q \right\}.$$

Lemma 3.14. Let $\{N_1, \ldots, N_q\}$ be as in (3.16), span $\{N_1, \ldots, N_q\} \cap \mathbb{S}_+^n = \{0\}$, and let

$$Q := \arg\max\{\log\det(X) : X = I + \sum_{i=1}^{q} x_i N_i > 0, \ x \in \mathbb{R}^q\}.$$
 (3.17)

Then for all $\alpha > 0$,

$$\alpha Q = \arg\max\{\log\det(X) : X = \alpha I + \sum_{i=1}^{q} x_i N_i \succ 0, \ x \in \mathbb{R}^q\}.$$
 (3.18)

Proof. We begin by expressing Q in terms of \mathcal{N} :

$$Q = \arg\max\{\log\det(X) : \mathcal{N}(X) = \mathcal{N}(I)\}.$$

By the assumption on the span of the matrices N_i and by Lemma 3.2, the feasible set of (3.17) is bounded. Moreover, the feasible set contains positive definite matrices, hence all the assumptions of Theorem 3.4 are satisfied. It follows that Q is the unique feasible, positive definite matrix satisfying $Q^{-1} \in \text{range}(\mathcal{N}^*)$.

Moreover, αQ is positive definite, feasible for (3.18), and $(\alpha Q)^{-1} \in \text{range}(\mathcal{N}^*)$. Therefore αQ is optimal for (3.18).

Now we prove that the parametric path converges to the analytic center under the condition of Lemma 3.14.

Theorem 3.15. Let $\{N_1, \ldots, N_q\}$ be as in (3.16). If $\operatorname{span}\{N_1, \ldots, N_q\} \cap \mathbb{S}^n_+ = \{0\}$ and \bar{X} is the limit point of the primal part of the parametric path as in Theorem 3.12, then $\bar{X} = \hat{X}$.

Proof. Let

$$\bar{X} =: \begin{bmatrix} \bar{Y} & 0 \\ 0 & 0 \end{bmatrix}, \ \hat{X} =: \begin{bmatrix} \hat{Y} & 0 \\ 0 & 0 \end{bmatrix}$$

and suppose, for eventual contradiction, that $\bar{Y} \neq \hat{Y}$. Then let $r, s \in \mathbb{R}$ be such that

$$\det(\bar{Y}) < r < s < \det(\hat{Y}).$$

Let Q be as in Lemma 3.14 and let $x \in \mathbb{R}^q$ satisfy $Q = I + \sum_{i=1}^q x_i N_i$. Now for any $\alpha > 0$ we have

$$\hat{X} + \alpha (I + \sum_{i=1}^{q} x_i A_i) = \begin{pmatrix} \hat{Y} + \alpha I + \alpha \sum_{i=1}^{q} x_i L_i & \alpha \sum_{i=1}^{q} x_i M_i \\ \alpha \sum_{i=1}^{q} x_i M_i^T & \alpha Q \end{pmatrix}.$$

Note that there exists $\varepsilon > 0$ such that $\hat{X} + \alpha \sum_{i=1}^{q} x_i A_i \succeq 0$ whenever $\alpha \in (0, \varepsilon)$. It follows that

$$\hat{X} + \alpha (I + \sum_{i=1}^{q} x_i A_i) \in \mathcal{F}(\alpha), \quad \forall \alpha \in (0, \varepsilon).$$

Taking the determinant, we have

$$\frac{1}{\alpha^{n-r}} \det(\hat{X} + \alpha(I + \sum_{i=1}^{q} x_i A_i)) = \frac{1}{\alpha^{n-r}} \det\left(\alpha Q - \alpha^2 (\sum_{i=1}^{q} x_i M_i)(\hat{Y} + \alpha I + \alpha \sum_{i=1}^{q} x_i L_i)^{-1} (\sum_{i=1}^{q} x_i M_i^T)\right) \\
\times \det(\hat{Y} + \alpha I + \alpha \sum_{i=1}^{q} x_i L_i),$$

$$= \det\left(Q - \alpha (\sum_{i=1}^{q} x_i M_i)(\hat{Y} + \alpha I + \alpha \sum_{i=1}^{q} x_i L_i)^{-1} (\sum_{i=1}^{q} x_i M_i^T)\right) \times \det(\hat{Y} + \alpha I + \alpha \sum_{i=1}^{q} x_i L_i).$$

Now we have

$$\lim_{\alpha \searrow 0} \frac{1}{\alpha^{n-r}} \det(\hat{X} + \alpha(I + \sum_{i=1}^{q} x_i A_i)) = \det(Q) \det(\hat{Y}).$$

Thus, there exists $\sigma \in (0, \varepsilon)$ so that for $\alpha \in (0, \sigma)$ we have

$$\det(\hat{X} + \alpha(I + \sum_{i=1}^{q} x_i A_i)) > s\alpha^{n-r} \det(Q).$$

As $X(\alpha)$ is the determinant maximizer over $\mathcal{F}(\alpha)$, we also have

$$\det(X(\alpha)) > s\alpha^{n-r} \det(Q), \quad \forall \alpha \in (0, \sigma). \tag{3.19}$$

On the other hand $X(\alpha) \to \bar{X}$ and let

$$X(\alpha) =: \begin{bmatrix} \alpha I + \sum_{i=1}^q x(\alpha)_i L_i & \sum_{i=1}^q x(\alpha)_i M_i \\ \sum_{i=1}^q x(\alpha)_i M_i^T & \alpha I + \sum_{i=1}^q x(\alpha)_i N_i \end{bmatrix}.$$

Then $\alpha I + \sum_{i=1}^{q} x(\alpha)_i L_i \to \bar{Y}$ and there exists $\delta \in (0, \sigma)$ such that for all $\alpha \in (0, \delta)$,

$$\det(\alpha I + \sum_{i=1}^{q} x(\alpha)_i L_i) < r.$$

Moreover, by definition of Q, $\det(\alpha I + \sum_{i=1}^{q} x(\alpha)_i N_i) \leq \det(\alpha Q) = \alpha^{n-r} \det(Q)$. To complete the proof, we apply the Hadamard-Fischer inequality to $\det(X(\alpha))$. For $\alpha \in (0, \delta)$ we have

$$\det(X(\alpha)) \le \det(\alpha I + \sum_{i=1}^{q} x(\alpha)_i L_i) \det(\alpha I + \sum_{i=1}^{q} x(\alpha)_i N_i) < r\alpha^{n-r} \det(Q),$$

a contradiction of (3.19).

Remark 3.16. Note that Example 3.13 fails the hypotheses of Theorem 3.15. Indeed, the matrix

$$\begin{bmatrix} 0 & 0 & 0 & 0 \\ 0 & 0 & 0 & -1 \\ 0 & 0 & 2 & 0 \\ 0 & -1 & 0 & 0 \end{bmatrix} lies in null(\mathcal{A}) and the bottom 2×2 block is nonzero and positive semidefinite.$$

4 The Projected Gauss-Newton Method

We have constructed a parametric path that converges to a point in the relative interior of \mathcal{F} . In this section we propose an algorithm to follow the path to its limit point. We do not prove convergence of the proposed algorithm and address its performance in Section 5. We follow the (projected) Gauss-Newton approach (the nonlinear analog of the Newton method) originally introduced for **SDPs** in [18] and improved more recently in [8]. This approach has been shown to have improved robustness compared to other symmetrization approaches. For well posed problems, the Jacobian for the search direction remains full rank in the limit to the optimum.

4.1 Scaled Optimality Conditions

The idea behind this approach is to view the system defining the parametric path as an overdetermined map and use the Gauss-Newton (GN) method for nonlinear systems. In the process, the linear feasibility equations are eliminated and the GN method is applied to the remaining bilinear equation. For $\alpha \geq 0$ let $G_{\alpha} : \mathbb{S}^{n}_{+} \times \mathbb{R}^{m} \times \mathbb{S}^{n}_{+} \to \mathbb{S}^{n} \times \mathbb{R}^{m} \times \mathbb{R}^{n \times n}$ be defined as

$$G_{\alpha}(X, y, Z) := \begin{bmatrix} \mathcal{A}^*(y) - Z \\ \mathcal{A}(X) - b(\alpha) \\ ZX - \alpha I \end{bmatrix}. \tag{4.1}$$

The solution to $G_{\alpha}(X, y, Z) = 0$ is exactly $(X(\alpha), y(\alpha), Z(\alpha))$ when $\alpha > 0$; and for $\alpha = 0$ the solution set is

$$\mathcal{F} \times (\mathcal{A}^*)^{-1}(\mathcal{D}) \times \mathcal{D}, \quad \mathcal{D} := \operatorname{range}(\mathcal{A}^*) \cap \operatorname{face}(\mathcal{F})^c.$$

Clearly, the limit point of the parametric path satisfies $G_0(X, y, Z) = 0$. We fix $\alpha > 0$. The GN direction, (dX, dy, dZ), uses the overdetermined GN system

$$G'_{\alpha}(X, y, Z) \begin{bmatrix} dX \\ dy \\ dZ \end{bmatrix} = -G_{\alpha}(X, y, Z).$$
 (4.2)

Note that the search direction is a strict descent direction for the norm of the residual, $\|\operatorname{vec}(G_{\alpha}(X,y,Z))\|_2^2$, when the Jacobian is full rank. The size of the problem is then reduced by projecting out the first two equations. We are left with a single linearization of the bilinear complementarity equation, i.e., n^2 equations in only t(n) variables. The least squares solution yields the projected GN direction after backsolves. We prefer steps of length 1, however, the primal and dual step lengths, α_p and α_d respectively, are reduced, when necessary, to ensure strict feasibility: $X + \alpha_p dX > 0$ and $Z + \alpha_d dZ > 0$. The parameter α is then reduced and the procedure repeated. On the parametric path, α satisfies

$$\alpha = \frac{\langle Z(\alpha), X(\alpha) \rangle}{n}.$$
(4.3)

Therefore, this is a good estimate of the target for α near the parametric path. As is customary, we then use a fixed $\sigma \in (0,1)$ to move the target towards optimality, $\alpha \leftarrow \sigma \alpha$.

4.1.1 Linearization and GN Search Direction

For the purposes of this discussion we vectorize the variables and data in G_{α} . Let $A \in \mathbb{R}^{m \times t(n)}$ be the matrix representation of A, that is

$$A_{i,:} := \operatorname{svec}(S_i)^T, \quad i \in \{1, \dots, m\}.$$

Let $N \in \mathbb{R}^{t(n) \times (t(n)-m)}$ be such that its columns form a basis for null(A) and let \hat{x} be a particular solution to $Ax = b(\alpha)$, e.g., the least squares solution. Then the affine manifold determined from the equation $\mathcal{A}(X) = b(\alpha)$ is equivalent to that obtained from the equation

$$x = \hat{x} + Nv, \quad v \in \mathbb{R}^{t(n)-m}.$$

Moreover, if $z := \operatorname{svec}(Z)$, we have the vectorization

$$g_{\alpha}(x, v, y, z) := \begin{bmatrix} A^{T}y - z \\ x - \hat{x} - Nv \\ sMat(z) sMat(x) - \alpha I \end{bmatrix} =: \begin{bmatrix} r_{d} \\ r_{p} \\ R_{c} \end{bmatrix}, \tag{4.4}$$

Now we show how the first two equations of the above system may be projected out, thereby reducing the size of the problem. First we have

$$g'_{\alpha}(x, v, y, z) \begin{pmatrix} dx \\ dv \\ dy \\ dz \end{pmatrix} = \begin{bmatrix} A^T dy - dz \\ dx - N dv \\ sMat(dz) sMat(x) + sMat(z) sMat(dx) \end{bmatrix},$$

and it follows that the GN step as in (4.2) is the least squares solution of the system

$$\begin{bmatrix} A^T dy - dz \\ dx - N dv \\ \operatorname{sMat}(dz) \operatorname{sMat}(x) + \operatorname{sMat}(z) \operatorname{sMat}(dx) \end{bmatrix} = - \begin{bmatrix} r_d \\ r_p \\ R_c \end{bmatrix}.$$

Since the first two equations are linear, we get $dz = A^T dy + r_d$ and $dx = N dv - r_p$. Substituting into the third equation we have,

$$\operatorname{sMat}(A^T dy + r_d) \operatorname{sMat}(x) + \operatorname{sMat}(z) \operatorname{sMat}(N dv - r_p) = -R_c.$$

After moving all the constants to the right hand side we obtain the projected GN system in dy and dv,

$$\operatorname{sMat}(A^T dy) \operatorname{sMat}(x) + \operatorname{sMat}(z) \operatorname{sMat}(N dv) = -R_c + \operatorname{sMat}(z) \operatorname{sMat}(r_p) - \operatorname{sMat}(r_d) \operatorname{sMat}(x). \tag{4.5}$$

The least squares solution to this system is the exact GN direction when $r_d = 0$ and $r_p = 0$, otherwise it is an approximation. We then use the equations $dz = A^T dy + r_d$ and $dx = N dv - r_p$ to obtain search directions for x and z.

In [8, Theorem 1], it is proved that if the solution set of $G_0(X, y, Z) = 0$ is a singleton such that X + Z > 0 and the starting point of the projected GN algorithm is sufficiently close to the parametric path then the algorithm, with a crossover modification, converges quadratically. As we showed above, the solution set to our problem is

$$\mathcal{F}\times (\mathcal{A}^*)^{-1}(\mathcal{D})\times \mathcal{D},$$

which is not a singleton as long as $\mathcal{F} \neq \emptyset$. Indeed, \mathcal{D} is a non-empty cone. Although the convergence result of [8] does not apply to our problem, their numerical tests indicate that the algorithm converges even for problems violating the strict complementarity and uniqueness assumptions and our observations agree.

4.2 Implementation Details

Several specific implementation modifications are used. We begin with initial x, v, y, z with corresponding $X, Z \succ 0$. If we obtain $P \succ 0$ as in Proposition 3.5 then we set Z = P and define y accordingly, otherwise Z = X = I. We estimate α using (4.3) and set $\alpha \leftarrow 2\alpha$ to ensure that our target is somewhat well centered to start.

4.2.1 Step Lengths and Linear Feasibility

We start with initial step lengths $\alpha_p = \alpha_d = 1.1$ and then backtrack using a Cholesky factorization test to ensure positive definiteness

$$X + \alpha_p dX > 0$$
, $Z + \alpha_d dZ > 0$.

If the step length we find is still > 1 after the backtrack, we set it to 1 and first update v, y and then update x, z using

$$x = \hat{x} + Nv, \quad z = A^T y.$$

This ensures exact linear feasibility. Thus we find that we maintain exact dual feasibility after a few iterations. Primal feasibility changes since α decreases. We have experimented with including an extra few iterations at the end of the algorithm with a fixed α to obtain exact primal feasibility (for the given α). In most cases the improvement of feasibility with respect to \mathcal{F} was minimal and not worth the extra computational cost.

4.2.2 Updating α and Expected Number of Iterations

In order to drive α down to zero, we fix $\sigma \in (0,1)$ and update alpha as $\alpha \leftarrow \sigma \alpha$. We use a moderate $\sigma = .6$. However, if this reduction is performed too quickly then our step lengths end up being too small and we get too close to the positive semidefinite boundary. Therefore, we change α using information from $\min\{\alpha_p, \alpha_d\}$. If the steplength is reasonably near 1 then we decrease using σ ; if the steplength is around .5 then we leave α as is; if the steplength is small then we increase to 1.2α ; and if the steplength is tiny (< .1), we increase to 2α . For most of the test problems, this strategy resulted in steplengths of 1 after the first few iterations.

We noted empirically that the condition number of the Jacobian for the least squares problem increases quickly, i.e., several singular values converge to zero. Despite this we are able to obtain high accuracy search directions.¹

Since we typically have steplengths of 1, α is generally decreased using σ . Therefore, for a desired tolerance ϵ and a starting $\alpha=1$ we would want $\sigma^k<\epsilon$, or equivalently, $k<\log_{10}(\epsilon)/\log_{10}(\sigma)$. For our $\sigma=.6$ and t decimals of desired accuracy, we expect to need k<4.5t iterations.

5 Generating Instances and Numerical Results

In this section we analyze the performance of an implementation of our algorithm. We begin with a discussion on generating spectrahedra. A particular challenge is in creating spectrahedra with specified singularity degree. Following this discussion, we present and analyze the numerical results.

¹Our algorithm finds the search direction using (4.5). If we looked at a singular value decomposition then we get the equivalent system $\Sigma(V^T d\bar{s}) = (U^T R H S)$. We observed that several singular values in Σ converge to zero while the corresponding elements in $(U^T R H S)$ converge to zero at a similar rate. This accounts for the improved accuracy despite the huge condition numbers. This appears to be a similar phenomenon to that observed in the analysis of interior point methods in [38,39] and as discussed in [13].

5.1 Generating Instances with Varying Singularity Degree

Our method for generating instances is motivated by the approach of [36] for generating **SDPs** with varying complementarity gaps. We begin by proving a relationship between strict complementarity of a primal-dual pair of **SDP** problems and the singularity degree of the optimal set of the primal **SDP**. This relationship allows us to modify the code presented in [36] and obtain spectrahedra having various singularity degrees. Recall the primal and dual **SDPs** in (2.5) and (2.6), respectively. Let $O_P \subseteq \mathbb{S}^n_+$ and $O_D \subseteq \mathbb{S}^n_+$ denote the primal and dual optimal sets respectively, where the dual optimal set is with respect to the variable Z. Specifically,

$$O_P := \{ X \in \mathbb{S}^n_+ : \mathcal{A}(X) = b, \ \langle C, X \rangle = p^* \}, \ O_D := \{ Z \in \mathbb{S}^n_+ : Z = C - \mathcal{A}^*(y), \ b^T y = d^*, \ y \in \mathbb{R}^m \}.$$

Note that O_P is a spectrahedron determined by the affine manifold

$$\begin{bmatrix} \mathcal{A}(X) \\ \langle C, X \rangle \end{bmatrix} = \begin{pmatrix} b \\ p^* \end{pmatrix}.$$

We note that the second system in the theorem of the alternative, Theorem 2.4, for the spectrahedron O_P is

$$0 \neq \tau C + \mathcal{A}^*(y) \succeq 0, \ \tau p^* + y^T b = 0.$$
 (5.1)

We say that strict complementarity holds for **SDP** and **D-SDP** if there exists $X^* \in O_P$ and $Z^* \in O_D$ such that

$$\langle X^{\star}, Z^{\star} \rangle = 0$$
 and $\operatorname{rank}(X^{\star}) + \operatorname{rank}(Z^{\star}) = n$.

If strict complementarity does not hold for **SDP** and **D-SDP** and there exist $X^* \in \operatorname{relint}(O_P)$ and $Z^* \in \operatorname{relint}(O_D)$, then we define the complementarity gap as

$$g := n - \operatorname{rank}(X^*) - \operatorname{rank}(Z^*).$$

Now we describe the relationship between strict complementarity of **SDP** and **D-SDP** and the singularity degree of O_P .

Proposition 5.1. If strict complementarity holds for SDP and D-SDP, then $sd(O_P) < 1$.

Proof. Let $X^* \in \operatorname{relint}(O_P)$. If $X^* \succ 0$, then $\operatorname{sd}(O_P) = 0$ and we are done. Thus we may assume $\operatorname{rank}(X^*) < n$. By strict complementarity, there exists $(y^*, Z^*) \in \mathbb{R}^m \times \mathbb{S}^n_+$ feasible for **D-SDP** with $Z^* \in O_D$ and $\operatorname{rank}(X^*) + \operatorname{rank}(Z^*) = n$. Now we show that $(1, -y^*)$ satisfies (5.1). Indeed, by dual feasibility,

$$C - \mathcal{A}^*(y^*) = Z^* \in \mathbb{S}^n_+ \setminus \{0\},\$$

and by complementary slackness,

$$p^* - (y^*)^T b = \langle X^*, C \rangle - \langle \mathcal{A}^*(y^*), X^* \rangle = \langle X^*, Z^* \rangle = 0.$$

Finally, since $\operatorname{rank}(X^*) + \operatorname{rank}(Z^*) = n$ we have $\operatorname{sd}(O_P) = 1$, as desired.

From the perspective of facial reduction, the interesting spectrahedra are those with singularity degree greater than zero and the above proposition gives us a way to construct spectrahedra with singularity degree exactly one. Using the algorithm of [36] we construct strictly complementary **SDPs** and then use the optimal set of the primal to construct a spectrahedron with singularity degree exactly one. Specifically, given positive integers n, m, r, and g the algorithm of [36] returns the data A, b, C corresponding to a primal dual pair of **SDPs**, together with $X^* \in \text{relint}(O_P)$ and $Z^* \in \text{relint}(O_D)$ satisfying

$$\operatorname{rank}(X^*) = r$$
, $\operatorname{rank}(Z^*) = n - r - g$.

Now if we set

$$\hat{\mathcal{A}}(X) := \begin{pmatrix} \mathcal{A}(X) \\ \langle C, X \rangle \end{pmatrix}, \ \hat{b} = \begin{pmatrix} b \\ \langle C, X^{\star} \rangle \end{pmatrix},$$

then $O_P = \mathcal{F}(\hat{\mathcal{A}}, \hat{b})$. Moreover, if g = 0 then $\mathrm{sd}(O_P) = 1$, by Proposition 5.1. This approach could also be used to create spectrahedra with larger singularity degrees by constructing **SDPs** with greater complementarity gaps, if the converse of Proposition 5.1 were true. We provide a sufficient condition for the converse in the following proposition.

Proposition 5.2. If $sd(O_P) = 0$, then strict complementarity holds for SDP and D-SDP. Moreover, if $sd(O_P) = 1$ and the set of solutions to (5.1) intersects $\mathbb{R}_{++} \times \mathbb{R}^m$, then strict complementarity holds for SDP and D-SDP.

Proof. Since we have only defined singularity degree for non-empty spectrahedra, there exists $X^* \in \operatorname{relint}(O_P)$. For the first statement, by Theorem 2.3, there exists $Z^* \in O_D$. Complementary slackness always holds, hence $\langle Z^*, X^* \rangle = 0$ and since $X^* \succ 0$ we have $Z^* = 0$. It follows that $\operatorname{rank}(X^*) + \operatorname{rank}(Z^*) = n$ and strict complementarity holds for **SDP** and **D-SDP**.

For the second statement, let $(\bar{\tau}, \bar{y})$ and $(\tilde{\tau}, \tilde{y})$ be solutions to (5.1) with $\bar{\tau} > 0$ and $\tilde{\tau}C + \mathcal{A}^*(\tilde{y})$ of maximal rank. Let

$$\bar{Z} := \bar{\tau}C + \mathcal{A}^*(\bar{\tau}), \ \tilde{Z} := \tilde{\tau}C + \mathcal{A}^*(\tilde{y}).$$

Then there exists $\varepsilon > 0$ such that $\bar{\tau} + \varepsilon \tilde{\tau} > 0$ and $\operatorname{rank}(\bar{Z} + \varepsilon \tilde{Z}) \ge \operatorname{rank}(\tilde{Z})$. Define

$$\tau := \bar{\tau} + \varepsilon \tilde{\tau}, \ y := \bar{y} + \varepsilon \tilde{y}, \ Z := \bar{Z} + \varepsilon \tilde{Z}.$$

Now (τ, y) is a solution to (5.1), i.e., $0 \neq \tau C + \mathcal{A}^*(y) \succeq 0$, $\tau p^* + y^T b = 0$. Moreover, $\operatorname{rank}(X^*) + \operatorname{rank}(Z) = n$ since $\operatorname{sd}(O_P) = 1$ and Z is of maximal rank. Now we define

$$Z^* := \frac{1}{\tau}Z = C - \mathcal{A}^* \left(-\frac{1}{\tau}y\right).$$

Since $\tau > 0$, it is clear that $Z^* \succeq 0$ and it follows that $\left(-\frac{1}{\tau}y, Z^*\right)$ is feasible for **D-SDP**. Moreover, this point is optimal since $d^* \geq -\frac{1}{\tau}y^Tb = p^* \geq d^*$. Therefore $Z^* \in O_D$ and since $\operatorname{rank}(Z^*) = \operatorname{rank}(Z)$, strict complementarity holds for **SDP** and **D-SDP**.

5.2 Numerical Results

n	m	r	$\lambda_1(X)$	$\lambda_r(X)$	$\lambda_{r+1}(X)$	$\lambda_n(X)$	$\ \mathcal{A}(X) - b\ _2$	$\langle Z, X \rangle$	α_f
50	100	25	1.06e+02	2.80e + 01	1.97e-11	5.07e-13	3.17e-12	1.26e-13	1.10e-12
80	160	40	8.74e + 01	3.22e+01	1.20e-10	9.00e-13	7.28e-12	2.95e-13	2.01e-12
110	220	55	7.74e + 01	3.73e+01	3.56e-10	7.23e-13	9.12e-12	3.65e-13	2.14e-12
140	280	70	7.82e + 01	3.84e + 01	4.11e-10	7.08e-13	1.26e-11	5.20e-13	2.65e-12

Table 5.1: Results for the case sd = 1. The eigenvalues refer to those of the primal variable, X, and each entry is the average of five runs.

For the numerical tests, we generate instances with $n \in \{50, 80, 110, 140\}$ and m = 2n. These are problems of small size relative to state of the art capabilities, nonetheless, we are able to demonstrate the performance of our algorithm through them. In Table 5.1 and Table 5.2 we record the results for the case sd = 1. For each instance, specified by n, m, and r, the results are the average of five runs. By r, we denote the maximum rank over all elements of the generated

n	m	r	$\lambda_1(Z)$	$\lambda_{r_d}(Z)$	$\lambda_{r_d+1}(Z)$	$\lambda_n(Z)$
50	100	25	1.85e+00	9.07e-02	3.96e-14	1.27e-14
80	160	40	1.96e+00	6.91e-02	6.23e-14	2.30e-14
110	220	55	1.98e+00	2.61e-02	5.77e-14	2.78e-14
140	280	70	2.03e+00	2.46e-02	6.96e-14	3.39e-14

Table 5.2: Eigenvalues of the dual variable, Z, corresponding to the primal variable of Table 5.1. Each entry is the average of five runs.

n	m	r	g	$\lambda_1(X)$	$\lambda_r(X)$	$\lambda_{r+1}(X)$	$\lambda_{r+g}(X)$	$\lambda_{r+g+1}(X)$	$\lambda_n(X)$	$ A(X) - b _2$	$\langle Z, X \rangle$	α_f
50	100	17	5	9.89e + 01	1.85e + 01	6.62e-05	2.61e-05	2.13e-10	6.10e-13	4.99e-12	2.04e-13	1.07e-12
80	160	27	8	1.11e+02	2.00e+01	1.89e-05	1.28e-05	7.36e-11	5.17e-13	8.40e-12	2.73e-13	1.27e-12
110	220	37	11	1.09e+02	2.42e+01	3.52e-05	2.33e-05	2.05e-10	1.52e-12	1.92e-11	6.46e-13	2.33e-12
140	280	47	14	1.63e+02	2.64e+01	1.07e-04	2.65e-05	1.02e-10	1.17e-13	9.84e-12	3.52e-13	1.48e-12

Table 5.3: Results for the case sd = 2. The eigenvalues refer to those of the primal variable, X, and each entry is the average of five runs.

n	m	r	g	$\lambda_1(Z)$	$\lambda_{r_d}(Z)$	$\lambda_{r_d+1}(Z)$	$\lambda_{r_d+g}(Z)$	$\lambda_{r_d+g+1}(Z)$	$\lambda_n(Z)$
50	100	17	5	2.22e+00	2.51e-02	1.04e-07	8.38e-08	9.18e-14	1.51e-14
80	160	27	8	2.03e+00	3.65e-02	1.03e-07	7.45e-08	7.92e-14	1.69e-14
110	220	37	11	2.13e+00	6.11e-02	1.78e-07	1.23e-07	1.36e-13	2.76e-14
140	280	47	14	2.19e+00	4.16e-02	7.39e-08	4.35e-08	6.04e-14	8.14e-15

Table 5.4: Eigenvalues of the dual variable, Z, corresponding to the primal variable of Table 5.3. Each entry is the average of five runs.

spectrahedron, which is fixed to r = n/2. In Table 5.1 we record the relevant eigenvalues of the primal variable, primal feasibility, complementarity, and the value of α at termination, denoted α_f . The values for primal feasibility and complementarity are sufficiently small and it is clear from the eigenvalues presented, that the first r eigenvalues are significantly smaller than the last n-r. These results demonstrate that the algorithm returns a matrix which is very close to the relative interior of \mathcal{F} . In Table 5.2 we record the relevant eigenvalues for the corresponding dual variable, Z. Note that $r_d := n - r$ and the eigenvalues recorded in the table indicate that Z is indeed an exposing vector. Moreover, it is a maximal rank exposing vector. While, we have not proved this, we observed that it is true for every test we ran with sd = 1.

In Table 5.3 and Table 5.4 we record similar values for problems where the singularity degree may be greater than 1. Using the approach described in Section 5.1 we generate instances of \mathbf{SDP} and $\mathbf{D}\text{-}\mathbf{SDP}$ having a complementarity gap of g and then we construct our spectrahedron from the optimal set of \mathbf{SDP} . By Proposition 5.1 and Proposition 5.2 the resulting spectrahedron may have singularity degree greater than 1. We observe that primal feasibility and complementarity are attained to a similar accuracy as in the $\mathrm{sd}=1$ case. The eigenvalues of the primal variable fall into three categories. The first r eigenvalues are sufficiently large so as not to be confused with 0, the last n-r-g eigenvalues are convincingly small, and the third group of eigenvalues, exactly g of them, are such that it is difficult to decide if they should be 0 or not. A similar phenomenon is observed for the eigenvalues of the dual variable. This demonstrates that exactly g of the eigenvalues are converging to 0 at a rate significantly smaller than that of the other n-r-g eigenvalues.

6 Conclusion

In this paper we have considered a *primal* approach to facial reduction for **SDPs** that reduces to finding a relative interior point of a spectrahedron. By considering a parametric optimization problem, we constructed a smooth path and proved that its limit point is in the relative interior of the spectrahedron. Moreover, we gave a sufficient condition for the relative interior point to coincide with the analytic center. We proposed a projected Gauss-Newton algorithm to follow the parametric path to the limit point and in the numerical results we observed that the algorithm converges. We also presented a method for constructing spectrahedra with singularity degree 1 and provided a sufficient condition for constructing spectrahedra of larger singularity degree. Finally, we showed that the parametric path has interesting structure for the simple cycle completion problem.

This research has also highlighted some new problems to be pursued. We single out two such problems. The first regards the eigenvalues of the limit point that are neither sufficiently small to be deemed zero nor sufficiently large to be considered as non-zero. We have experimented with some eigenvalue deflation techniques, but none have led to a satisfactory method. Secondly, there does not seem to be a method in the literature for constructing spectrahedra with specified singularity degree.

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