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**On the Norms Used in Computing
the Structured Singular Value**

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Abstract

Different norms are considered to replace the Euclidean norm in an algorithm given by Fan and Tits (IEEE Trans. Automat. Contr., AC-33, pp.284-289, 1988) which is used for the computation of the structured singular value of any matrix. It is shown that the ℓ_1 -norm is the best possible norm in a certain sense.

Recently, there has been a considerable amount of interest in the study of the structured singular value, the concept of which was originated by Doyle [1] and is used as a tool for the analysis and synthesis of feedback systems with structured uncertainties (e.g., see [2,3] and their references).

Let M be an $n \times n$ complex matrix, and $\mathcal{K} = (k_1, \dots, k_m)$ an m -tuple of positive integers which satisfies $\sum_{i=1}^m k_i = n$. For $i = 1, \dots, m$, denote the i th-block-projection matrix by $P_i = \text{block diag}(O_{k_1}, \dots, O_{k_{i-1}}, I_{k_i}, O_{k_{i+1}}, \dots, O_{k_m})$, where O_k and I_k are the zero matrix and identity matrix, respectively, of order $k \times k$ for any positive integer k . Then the structured singular value of M with respect to the block-structure \mathcal{K} is the nonnegative scalar

$$\mu(M) = \max_{x \in \partial B} \{ \|Mx\| : \|P_i x\| \|Mx\| = \|P_i Mx\|, i = 1, \dots, m \},$$

where $\|\cdot\|$ denotes the Euclidean (ℓ_2 -) norm in \mathbb{C}^n and ∂B the corresponding unit sphere. One major issue in the study of $\mu(M)$ is the computation of it. In [2], the authors devise an algorithm [2, Algorithm 1], which we shall explain immediately, to compute $\mu(M)$. They first define, for any real number α , the hermitian matrices

$$A_i(\alpha) = \alpha P_i - M^H P_i M, \text{ for } i = 1, \dots, m,$$

and the m -form numerical range associated with $A_1(\alpha), \dots, A_m(\alpha)$

$$W(\alpha) = \{ (v_1, \dots, v_m) \in \mathbb{R}^m : \exists x \in \partial B \text{ such that } v_i = x^H A_i(\alpha) x \text{ for all } i \}.$$

A function $c(\cdot) : \mathbb{R} \rightarrow \mathbb{R}$, which depends on M and \mathcal{K} , is then defined by

$$c(\alpha) = \min \{ \|v\| : v \in W(\alpha) \},$$

where $\|\cdot\|$ denotes again the Euclidean norm (on \mathbb{R}^m this time). Then they show [2, Corollary 1 and Proposition 1] that, for any matrix M with structure \mathcal{K} , $c(\cdot)$ satisfies

$$c(\cdot) \text{ is continuous,} \tag{1}$$

$$c(\mu^2) = 0, \ c(\alpha) > 0 \text{ for all } \alpha > \mu^2, \tag{2}$$

and

$$c(\alpha + s) \leq c(\alpha) + s \text{ for all } s \geq 0 \text{ and real } \alpha, \quad (3)$$

where $\mu = \mu(M)$. The initial step in Algorithm 1 of [2] is to set $\alpha_0 = \bar{\sigma}^2(M)$, where $\bar{\sigma}(M)$ is the largest singular value of M . The iteration step is to set $\alpha_{k+1} = \alpha_k - c(\alpha_k)$ for $k = 0, 1, 2, \dots$. The authors show in [2, Theorem 2] that, since $c(\cdot)$ satisfies (1) through (3), Algorithm 1 will generate a monotonic decreasing sequence $\{\alpha_k\}$ with $\lim_{k \rightarrow \infty} \alpha_k = \mu^2$. The structured singular value μ can thus be obtained. They also remark in the footnote that the Euclidean norm in the definition of $c(\cdot)$ can be replaced by the ℓ_1 -norm to get the strongest version of Proposition 1. The purpose of the present note is to elaborate this remark.

Suppose we consider any norm $N(\cdot)$ on \mathbb{R}^m instead of the Euclidean norm. Similar to the case of $c(\cdot)$, we may define a function $c_N : \mathbb{R} \rightarrow \mathbb{R}$ (which also depends on M and \mathcal{K}) by

$$c_N(\alpha) = \min\{N(v) : v \in W(\alpha)\}.$$

It is not hard to see that $c_N(\cdot)$ always satisfies (1) and (2); and if in addition $c_N(\cdot)$ satisfies (3) then the function $c(\cdot)$ in Algorithm 1 can be replaced by $c_N(\cdot)$. Let

$$\mathcal{N} = \{N(\cdot) : N(\cdot) \text{ is a norm on } \mathbb{R}^m \text{ and } c_N(\cdot) \text{ satisfies} \\ \text{condition (3) for all matrices } M \text{ with structure } \mathcal{K} \}.$$

Then Algorithm 1 will work with $c(\cdot)$ being replaced by any $c_N(\cdot)$ where $N(\cdot) \in \mathcal{N}$. In view of the iteration step of the algorithm, we may want to choose a norm $N_0(\cdot) \in \mathcal{N}$ such that

$$c_{N_0}(\alpha) \geq c_N(\alpha) \text{ for all } N(\cdot) \in \mathcal{N} \text{ and } \alpha > \mu^2,$$

so that the resulting algorithm has the fastest convergent rate and is thus the most efficient. The following result shows that $\|\cdot\|_1$, i.e., the ℓ_1 -norm on \mathbb{R}^m defined by

$$\|v\|_1 = \sum_{i=1}^m |v_i| \text{ for all } v = (v_1, \dots, v_m) \in \mathbb{R}^m,$$

will give such a “best possible” norm.

Theorem. Let $\mathcal{K} = (k_1, \dots, k_m)$ be a given block structure. Then

- (a) $\|\cdot\|_1 \in \mathcal{N}$, and
- (b) for any $N(\cdot) \in \mathcal{N}$, $\|v\|_1 \geq N(v)$ for all $v \in \mathbb{R}^m$,

so that

$$c_1(\alpha) := \min\{\|v\|_1 : v \in W(\alpha)\} \geq c_N(\alpha)$$

for any real α and complex matrix M with block structure \mathcal{K} .

Proof. (a) The proof of the fact that $c_1(\cdot)$ satisfies (3) is similar to (and simpler than) that of $c(\cdot)$ given in the proof of Proposition 1 in [2]. Hence $\|\cdot\|_1 \in \mathcal{N}$.

(b) We prove by contradiction. Suppose $N(\cdot) \in \mathcal{N}$ but $N(v) > \|v\|_1$ for some $v \in \mathbb{R}^m$. Let $\{e_1, \dots, e_m\}$ be the standard basis for \mathbb{R}^m . Without loss of generality we may assume

$$N(e_1) = 1 + q > \|e_1\|_1 = 1;$$

for if $N(e_i) \leq \|e_i\|_1$ for all $i = 1, \dots, m$, then

$$N(v) = N\left(\sum_{i=1}^m v_i e_i\right) \leq \sum_{i=1}^m |v_i| N(e_i) \leq \sum_{i=1}^m |v_i| \|e_i\|_1 = \|v\|_1$$

for all $v = (v_1, \dots, v_m) \in \mathbb{R}^m$, which contradicts the assumption. Let

$$L = \{v \in \mathbb{R}^m : v_1 + w_2 v_2 + \dots + w_m v_m = 1\}$$

be a supporting hyperplane to the ball in \mathbb{R}^m with respect to $N(\cdot)$ with radius $(1 + q)$.

As e_1 is on L , we have

$$\min_{v \in L} N(v) = N(e_1) = 1 + q.$$

Note that some of the w_i 's may be zero. However, for any $\epsilon > 0$, we can choose nonzero $w'_2, \dots, w'_m \in \mathbb{R}$ which are arbitrary close to w_2, \dots, w_m respectively, such that if

$$L' = \{v \in \mathbb{R}^m : v_1 + w'_2 v_2 + \dots + w'_m v_m = 1\}$$

then

$$\min_{v \in L'} N(v) > N(e_1) - \epsilon.$$

Now choose $\lambda, \beta_1, \dots, \beta_m \in \mathbb{R}$ such that $\lambda - \beta_1^2 = 1$ and $\lambda - \beta_i^2 = 1/w'_i$ for $i = 2, \dots, m$. Then, by defining

$$M = \sum_{i=1}^m \beta_i P_i \tag{4}$$

(which depends on ϵ , because β_i 's depend on ϵ), we have, for any real α ,

$$\begin{aligned} A_i(\alpha) &= \alpha P_i - \left(\sum_{j=1}^m \beta_j P_j \right) P_i \left(\sum_{j=1}^m \beta_j P_j \right) \\ &= (\alpha - \beta_i^2) P_i \text{ for } i = 1, \dots, m \end{aligned}$$

and

$$\begin{aligned} c_N(\alpha) &= \min \{ N(v) : v_i = x^H A_i(\alpha) x, \ i = 1, \dots, m, \ x \in \partial B \} \\ &= \min \{ N(v) : v_i = (\alpha - \beta_i^2) a_i, \ a_i \geq 0, \ i = 1, \dots, m, \\ &\quad a_1 + \dots + a_m = 1 \}. \end{aligned}$$

Hence

$$\begin{aligned} c_N(\lambda - 1) &= \min \{ N(v) : v_i = (\lambda - 1 - \beta_i^2) a_i, \ a_i \geq 0, \ i = 1, \dots, m, \\ &\quad a_1 + \dots + a_m = 1 \} \\ &= \min \{ N(v) : v_1 = 0, \ v_i = (\lambda - 1 - \beta_i^2) a_i, \ a_i \geq 0, \\ &\quad i = 2, \dots, m, \ a_1 \geq 0, \ a_1 + \dots + a_m = 1 \} . \end{aligned}$$

By putting $(a_1, \dots, a_m) = e_1$, we get

$$c_N(\lambda - 1) = N(0) = 0.$$

Also,

$$\begin{aligned}
c_N(\lambda) &= \min \{N(v) : v_i = (\lambda - \beta_i^2)a_i, a_i \geq 0, i = 1, \dots, m, \\
&\quad a_1 + \dots + a_m = 1\} \\
&= \min \{N(v) : v_1 = a_1 \geq 0, v_i = a_i/w'_i, a_i \geq 0, i = 2, \dots, m, \\
&\quad a_1 + \dots + a_m = 1\} \\
&= \min \{N(v) : v_1 + w'_2 v_2 + \dots + w'_m v_m = 1, v_1 \geq 0, \\
&\quad w'_i v_i \geq 0, i = 2, \dots, m\} \\
&\geq \min \{N(v) : v \in L'\} \\
&> N(e_1) - \epsilon \\
&= (1 + q) - \epsilon \\
&= (1 + q) + c_N(\lambda - 1) - \epsilon.
\end{aligned}$$

As $\epsilon > 0$ is arbitrary, we may choose $0 < \epsilon < q$ so that, for the matrix M defined in (4), we have

$$c_N((\lambda - 1) + 1) = c_N(\lambda) > c_N(\lambda - 1) + 1.$$

As a result, $c_N(\cdot)$ does not satisfy (3) for this matrix M , and hence $N(\cdot) \notin \mathcal{N}$. Thus, if $N(\cdot) \in \mathcal{N}$ then we must have

$$\|v\|_1 \geq N(v) \text{ for all } v \in \mathbb{R}^m,$$

and hence

$$\begin{aligned}
c_1(\alpha) &= \min \{\|v\|_1 : v \in W(\alpha)\} \\
&\geq \min \{N(v) : v \in W(\alpha)\} \\
&= c_N(\alpha) \text{ for any real } \alpha.
\end{aligned}$$

□

Now the problem remains to devise a method for computing $c_1(\alpha)$, so that the algorithm can be implemented. The computation of $c_1(\alpha)$ (or $c(\alpha)$) for any given M

and \mathcal{K} can be, in general, difficult. However, similar to the case of $c(\alpha)$, there are existing methods for computing the value

$$c'_1(\alpha) := \min\{\|v\|_1 : v \in \text{co } W(\alpha)\},$$

where $\text{co } W(\alpha)$ denotes the convex hull of $W(\alpha)$. Let $\langle \cdot, \cdot \rangle$ denote the usual inner product in \mathbb{R}^m . Since the ℓ_∞ -norm ($\|\cdot\|_\infty$) is the dual norm of the ℓ_1 -norm, we have

$$\begin{aligned} c'_1(\alpha) &= \min_{v \in \text{co } W(\alpha)} \|v\|_1 \\ &= \min_{v \in \text{co } W(\alpha)} \max_{\substack{a \in \mathbb{R}^m \\ \|a\|_\infty \leq 1}} \langle v, a \rangle. \end{aligned} \tag{5}$$

As $\text{co } W(\alpha)$ and $\{a \in \mathbb{R}^m : \|a\|_\infty \leq 1\}$ are convex sets, and $\langle \cdot, \cdot \rangle$ is bilinear, the max and min in (5) can be interchanged to yield

$$\begin{aligned} c'_1(\alpha) &= \max_{\substack{a \in \mathbb{R}^m \\ \|a\|_\infty \leq 1}} \min_{v \in \text{co } W(\alpha)} \langle v, a \rangle \\ &= \max_{|a_i| \leq 1} \min_{x \in \partial B} \sum_{i=1}^m a_i x^H A_i(\alpha) x \\ &= \max_{|a_i| \leq 1} \min_{x \in \partial B} x^H \left(\sum_{i=1}^m a_i A_i(\alpha) \right) x \\ &= \max_{|a_i| \leq 1} \lambda_{\min} \left(\sum_{i=1}^m a_i A_i(\alpha) \right) \\ &= - \min_{|a_i| \leq 1} \lambda_{\max} \left(\sum_{i=1}^m a_i A_i(\alpha) \right), \end{aligned} \tag{6}$$

where λ_{\min} and λ_{\max} denote respectively the smallest and largest eigenvalue. Several existing algorithms are available for solving the convex problem (6). We refer to a recent paper by Boyd and Yang [4, Section 7] the discussion on the details, including the advantages and disadvantages, of these numerical algorithms. As a result, if $W(\alpha)$ is convex (which is always the case when $m \leq 3$; see [5] for example), $c_1(\alpha)$ ($= c'_1(\alpha)$) could be computed with any of these algorithms. If $m \geq 4$, then $W(\alpha)$ may not be convex. In

this case, replacing $c(\cdot)$ by $c'_1(\cdot)$ in Algorithm 1 of [2] will result in a sequence $\{\sqrt{\alpha_k}\}$ with limit

$$\mu' = \inf \{ \sqrt{\alpha} : c'_1(\beta) > 0 \text{ for all } \beta \geq \alpha \},$$

which is clearly an upper bound for $\mu(M)$. This is exactly the same situation as the case of $c(\cdot)$ and $c'(\cdot)$ discussed in [2].

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