

A Greedy Subspace Method for Computing the \mathcal{L}_∞ -Norm

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We consider the computation of the \mathcal{L}_∞ -norm for a general class of \mathcal{L}_∞ -functions and focus on the case where the function is represented in terms of large-scale matrix-valued factors. We propose a subspace projection method to obtain reduced approximations of this function by interpolation techniques. The \mathcal{L}_∞ -norms are computed for the resulting reduced functions, then the subspaces are refined by means of the optimizer of the \mathcal{L}_∞ -norm of the reduced function. In this way we obtain much better performance compared to existing methods.

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1 Problem Statement

In this paper we consider functions of the form

$$H : \Omega \rightarrow \mathbb{C}^{p \times m}, \quad H(s) := C(s)D(s)^{-1}B(s), \quad (1)$$

where $\Omega \subseteq \mathbb{C}$ is assumed to be an open domain enclosing the imaginary axis $i\mathbb{R}$. Moreover, we assume that the matrix-valued functions $B : \Omega \rightarrow \mathbb{C}^{n \times m}$, $C : \Omega \rightarrow \mathbb{C}^{p \times n}$, and $D : \Omega \rightarrow \mathbb{C}^{n \times n}$ are defined by

$$\begin{aligned} B(s) &:= f_1(s)B_1 + \dots + f_{\kappa_B}(s)B_{\kappa_B}, \\ C(s) &:= g_1(s)C_1 + \dots + g_{\kappa_C}(s)C_{\kappa_C}, \\ D(s) &:= h_1(s)D_1 + \dots + h_{\kappa_D}(s)D_{\kappa_D}, \end{aligned} \quad (2)$$

for given matrices $B_1, \dots, B_{\kappa_B} \in \mathbb{C}^{n \times m}$, $C_1, \dots, C_{\kappa_C} \in \mathbb{C}^{p \times n}$, $D_1, \dots, D_{\kappa_D} \in \mathbb{C}^{n \times n}$ and given functions f_1, \dots, f_{κ_B} , g_1, \dots, g_{κ_C} , $h_1, \dots, h_{\kappa_D} : \Omega \rightarrow \mathbb{C}$ that are meromorphic in Ω .

In this paper we suppose the function H (more precisely, its restriction to the imaginary axis) to be an element of the normed space

$$\mathcal{L}_\infty^{p \times m} := \left\{ H|_{i\mathbb{R}} \mid H : \Omega \rightarrow \mathbb{C}^{p \times m} \text{ is analytic for an open domain } \Omega \subseteq \mathbb{C} \text{ with } i\mathbb{R} \subset \Omega \text{ and } \sup_{\omega \in \mathbb{R}} \|H(i\omega)\|_2 < \infty \right\}.$$

For ease of notation, we write $H \in \mathcal{L}_\infty^{p \times m}$ instead of $H|_{i\mathbb{R}} \in \mathcal{L}_\infty^{p \times m}$. The \mathcal{L}_∞ -norm of a function $H \in \mathcal{L}_\infty^{p \times m}$ is defined by

$$\|H\|_{\mathcal{L}_\infty} := \sup_{\omega \in \mathbb{R}} \|H(i\omega)\|_2 = \sup_{\omega \in \mathbb{R}} \sigma_{\max}(H(i\omega)),$$

where $\sigma_{\max}(\cdot)$ denotes the maximum singular value of its matrix argument. In this paper we address the computation of the \mathcal{L}_∞ -norm with an emphasis on the case where n is large with $n \gg m, p$ and thus the evaluation of $H(s)$ is expensive.

2 The Method

Our approach makes use of ideas from the field of model order reduction, i. e., instead of computing the \mathcal{L}_∞ -norm of H directly, we construct reduced functions of the form

$$\tilde{H} : \Omega \rightarrow \mathbb{C}^{p \times m}, \quad \tilde{H}(s) := \tilde{C}(s)\tilde{D}(s)^{-1}\tilde{B}(s), \quad (3)$$

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where

$$\begin{aligned}\tilde{B}(s) &:= f_1(s)\tilde{B}_1 + \cdots + f_{\kappa_B}(s)\tilde{B}_{\kappa_B}, & \tilde{B}_j &:= W^*B_j, \quad j = 1, \dots, \kappa_B, \\ \tilde{C}(s) &:= g_1(s)\tilde{C}_1 + \cdots + g_{\kappa_C}(s)\tilde{C}_{\kappa_C}, & \tilde{C}_j &:= C_jV, \quad j = 1, \dots, \kappa_C, \\ \tilde{D}(s) &:= h_1(s)\tilde{D}_1 + \cdots + h_{\kappa_D}(s)\tilde{D}_{\kappa_D}, & \tilde{D}_j &:= W^*D_jV, \quad j = 1, \dots, \kappa_D,\end{aligned}$$

and the matrices $V, W \in \mathbb{C}^{n \times k}$ with $k \ll n$ are appropriately chosen. Since \tilde{D} is of much smaller dimension than D , the \mathcal{L}_∞ -norm of \tilde{H} can be computed much more efficiently.

In our method we compute a sequence of such reduced functions $\tilde{H}_1, \tilde{H}_2, \dots$, whose \mathcal{L}_∞ -norms converge to $\|H\|_{\mathcal{L}_\infty}$. For the construction of the reduced functions we use the interpolation technique of [1]. Assume first that $m = p$ and that r interpolation points $i\omega_1, \dots, i\omega_r$ are given. Then \tilde{H}_r is defined by \tilde{H} as in (3) with V, W set equal to the particular choices

$$V_r = [D(i\omega_1)^{-1}B(i\omega_1) \quad \dots \quad D(i\omega_r)^{-1}B(i\omega_r)], \quad W_r = [D(i\omega_1)^{-*}C(i\omega_1)^* \quad \dots \quad D(i\omega_r)^{-*}C(i\omega_r)^*]. \quad (5)$$

These choices for V and W give rise to the Hermite interpolation properties [1, 2]

$$H(i\omega_k) = \tilde{H}_r(i\omega_k), \quad H'(i\omega_k) = \tilde{H}'_r(i\omega_k), \quad k = 1, \dots, r. \quad (6)$$

Note that the number of columns of V_r and W_r have to be equal so that $\tilde{D}(s)^{-1}$ exists. This is violated by the choices in (5) unless $m = p$. The following choices overcome this difficulty when $m \neq p$ while preserving the Hermite interpolation properties (6):

$$\begin{aligned}V_r &= [D(i\omega_1)^{-1}B(i\omega_1)H(i\omega_1)^* \quad \dots \quad D(i\omega_r)^{-1}B(i\omega_r)H(i\omega_r)^*], & W_r &\text{ as in (5),} & \text{if } m > p, \\ W_r &= [D(i\omega_1)^{-*}C(i\omega_1)^*H(i\omega_1) \quad \dots \quad D(i\omega_r)^{-*}C(i\omega_r)^*H(i\omega_r)], & V_r &\text{ as in (5),} & \text{if } m < p.\end{aligned}$$

Once we have constructed \tilde{H}_r , we compute its \mathcal{L}_∞ -norm by established methods for the small-scale case such as [3–5]. Using the point $i\omega_{r+1}$ where the \mathcal{L}_∞ -norm of \tilde{H}_r is attained, we update the projection spaces

$$\begin{aligned}V_{r+1} &:= [V_r \quad D(i\omega_{r+1})^{-1}B(i\omega_{r+1})] \quad (\text{or } V_{r+1} := [V_r \quad D(i\omega_{r+1})^{-1}B(i\omega_{r+1})H(i\omega_{r+1})^*]), \\ W_{r+1} &:= [W_r \quad D(i\omega_{r+1})^{-*}C(i\omega_{r+1})^*] \quad (\text{or } W_{r+1} := [W_r \quad D(i\omega_{r+1})^{-*}C(i\omega_{r+1})^*H(i\omega_{r+1})]).\end{aligned}$$

Under the assumption of convergence (that we usually observe), we can prove a superlinear rate of convergence to a local maximizer of $\sigma_{\max}(H(i\cdot))$. The proof makes use of the Hermite interpolation property (6) and is elaborated in [2].

3 A Numerical Example

A MATLAB implementation of our method has been made available on the internet¹. In this section we briefly demonstrate this implementation on a numerical example. As a benchmark we take the `S80PI_n` example², which is a descriptor system with $n = 4182$ states and $m = p = 1$ input and output, respectively. We obtain a transfer function $H(s) = C(sE - A)^{-1}B$ which is of the form (1). We start initially with 10 equally spaced interpolation points in the interval $[0.1, 10000]$. Only 5 additional iterations are needed so that the points where the \mathcal{L}_∞ -norms of two consecutive reduced functions are attained are located at a relative distance of at most 10^{-6} . Our method correctly computes $\|H\|_{\mathcal{L}_\infty} = 3.37016$ which is attained at $\omega = 6961.49$. The runtime is only 0.20 s compared to 3.96 s required by the pseudospectral method from [6].

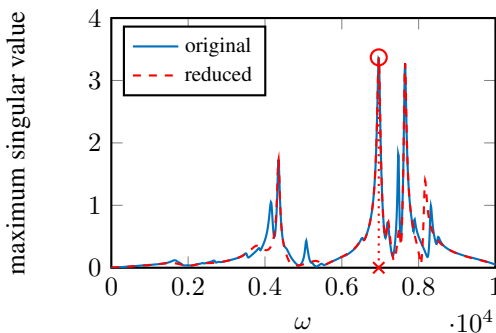


Fig. 1 Original and final reduced transfer functions obtained for the `S80PI_n` example. The red cross and circle indicate the location of the maximizer and the \mathcal{L}_∞ -norm of the reduced function, respectively.

¹ see <http://www.math.tu-berlin.de/index.php?id=186267&L=1>

² available from <http://sites.google.com/site/rommes/software>

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