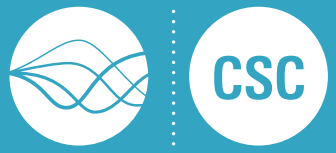
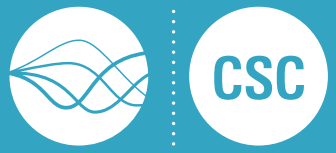

MAX PLANCK INSTITUTE
FOR DYNAMICS OF COMPLEX
TECHNICAL SYSTEMS
MAGDEBURG
CSC
COMPUTATIONAL METHODS IN
SYSTEMS AND CONTROL THEORY
Parametric Model Order Reduction for
Linear Control Systems
Peter Benner
HRZZ Project
"Control of Dynamical Systems (ConDys)"
— second project meeting —
Zagreb, 2–3 November 2017
ConDyS



Outline



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Introduction

Parametric Dynamical Systems

$$\Sigma(p): \begin{cases} E(p)\dot{x}(t;p) = f(t, x(t;p), u(t), p), \quad x(t_0) = x_0, & \text{(a)} \\ y(t;p) = g(t, x(t;p), u(t), p) & \text{(b)} \end{cases}$$

with

- (generalized) states $x(t;p) \in \mathbb{R}^n$ ($E \in \mathbb{R}^{n\times n}$),
- inputs (controls) $u(t) \in \mathbb{R}^m$,
- outputs (measurements, quantity of interest) $y(t;p) \in \mathbb{R}^q$,
  (b) is called output equation,
- $p \in \Omega \subset \mathbb{R}^d$ is a parameter vector, Ω is bounded.

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$E(p)$ singular $\Rightarrow$ (a) is system of differential-algebraic equations (DAEs)
otherwise $\Rightarrow$ (a) is system of ordinary differential equations (ODEs)

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Applications:

- Repeated simulation for varying material or geometry parameters, boundary conditions,
- control, optimization and design,
- of models, often generated by FE software (e.g., ANSYS, NASTRAN,...) or automatic tools (e.g., Modelica).

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Parametric discretized model often not available,
but matrices for certain parameter values can be extracted

(or output data for given u and p can be generated!)

Introduction

Linear, Time-Invariant (Parametric) Systems

$$\begin{aligned} E(p)\dot{x}(t;p) &= A(p)x(t;p) + B(p)u(t), & A(p),\ E(p) &\in \mathbb{R}^{n\times n}, \\ y(t;p) &= C(p)x(t;p), & B(p) &\in \mathbb{R}^{n\times m},\ C(p) \in \mathbb{R}^{q\times n}. \end{aligned}$$

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Laplace Transformation / Frequency Domain

Application of Laplace transformation

$$x(t;p) \mapsto x(s;p), \quad \dot{x}(t;p) \mapsto sx(s;p)$$

to linear system with $x(0;p) \equiv 0$:

$$sE(p)x(s;p) = A(p)x(s;p) + B(p)u(s), \quad y(s;p) = C(p)x(s;p),$$

yields I/O-relation in frequency domain:

$$y(s;p) = \Big(\underbrace{C(p)(sE(p) - A(p))^{-1}B(p)}_{=:G(s,p)} \Big) u(s).$$

$G(s,p)$ is the parameter-dependent transfer function of $\Sigma(p)$.

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Goal: Fast evaluation of mapping $(u,p) \ \to \ y(s;p)$.

The Parametric Model Order Reduction (PMOR) Problem

Problem

Approximate the dynamical system

$$\begin{aligned} E(p)\dot{x} &= A(p)x + B(p)u, & E(p),\ A(p) &\in \mathbb{R}^{n\times n}, \\ y &= C(p)x, & B(p) &\in \mathbb{R}^{n\times m},\ C(p) \in \mathbb{R}^{q\times n}, \end{aligned}$$

by reduced-order system

$$\begin{aligned} \hat{E}(p)\dot{\hat{x}} &= \hat{A}(p)\hat{x} + \hat{B}(p)u, & \hat{E}(p),\ \hat{A}(p) &\in \mathbb{R}^{r\times r}, \\ \hat{y} &= \hat{C}(p)\hat{x}, & \hat{B}(p) &\in \mathbb{R}^{r\times m},\ \hat{C}(p) \in \mathbb{R}^{q\times r}, \end{aligned}$$

of order $r \ll n$, such that

$$\|y - \hat{y}\| = \|Gu - \hat{G}u\| \leq \|G - \hat{G}\| \cdot \|u\| < \text{tolerance} \cdot \|u\| \quad \forall\ p \in \Omega.$$

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$\Longrightarrow$ Approximation problem: $\min\limits_{\text{order}\,(\hat{G}) \leq r} \|G - \hat{G}\|.$

The Parametric Model Order Reduction (PMOR) Problem

Structure Preservation

Parametric System

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Parametric model reduction goal:

preserve parameters as *symbolic quantities* in reduced-order model:

$$\widehat{\Sigma}(p): \begin{cases} \hat{E}(p)\dot{\hat{x}}(t;p) &= \hat{A}(p)\hat{x}(t;p) + \hat{B}(p)u(t), \\ \hat{y}(t;p) &= \hat{C}(p)\hat{x}(t;p) \end{cases}$$

with states $\hat{x}(t;p) \in \mathbb{R}^r$ and $r \ll n$.

The Parametric Model Order Reduction (PMOR) Problem

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Assuming parameter-affine representation:

$$\begin{aligned} E(p) &= E_0 + e_1(p)E_1 + \ldots + e_{q_E}(p)E_{q_E}, \\ A(p) &= A_0 + a_1(p)A_1 + \ldots + a_{q_A}(p)A_{q_A}, \\ B(p) &= B_0 + b_1(p)B_1 + \ldots + b_{q_B}(p)B_{q_B}, \\ C(p) &= C_0 + c_1(p)C_1 + \ldots + c_{q_C}(p)C_{q_C}, \end{aligned}$$

allows easy parameter preservation for projection based model reduction.

Structure Preservation

Petrov-Galerkin-type projection

For given projection matrices $V, W \in \mathbb{R}^{n\times r}$ with $W^T V = I_r$ ($\rightsquigarrow (VW^T)^2 = VW^T$ is projector), compute

$$\begin{aligned}
\hat{E}(p) &= W^T E_0 V + e_1(p) W^T E_1 V + \ldots + e_{q_E}(p) W^T E_{q_E} V \\
\hat{A}(p) &= W^T A_0 V + a_1(p) W^T A_1 V + \ldots + a_{q_A}(p) W^T A_{q_A} V \\
\hat{B}(p) &= W^T B_0 \quad + b_1(p) W^T B_1 \quad + \ldots + b_{q_B}(p) W^T B_{q_B} \\
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\hat{A}(p) &= W^T A_0 V + a_1(p) W^T A_1 V + \ldots + a_{q_A}(p) W^T A_{q_A} V \\
&= \hat{A}_0 + a_1(p)\hat{A}_1 + \ldots + a_{q_A}(p)\hat{A}_{q_A} \\
\hat{B}(p) &= W^T B_0 \quad + b_1(p) W^T B_1 \quad + \ldots + b_{q_B}(p) W^T B_{q_B} \\
&= \hat{B}_0 + b_1(p)\hat{B}_1 + \ldots + b_{q_B}(p)\hat{B}_{q_B} \\
\hat{C}(p) &= \quad C_0 V + \quad c_1(p) C_1 V + \ldots + \quad c_{q_C}(p) C_{q_C} V \\
&= \hat{C}_0 + c_1(p)\hat{C}_1 + \ldots + c_{q_C}(p)\hat{C}_{q_C}
\end{aligned}$$

Basis Generation — Global vs. Local

Local Bases

Obtain $V_k, W_k \in \mathbb{R}^{n \times r_k}$ using any non-parametric linear MOR method for a number of full-order models $\Sigma(p^{(k)})$, $k = 1, \ldots, \ell$. Then compute reduced-order model by

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no need for affine parametrization, requires only system matrices $A(p^{(k)}), B(p^{(k)}), \ldots$.

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2. transfer function interpolation: spurious poles of the parametric transfer function.
3. matrix interpolation: different models obtained in different coordinate systems $\rightsquigarrow$ Procrustes problem $\rightsquigarrow$ potential loss of accuracy; efficiency in "online" phase suffers from evaluating the interpolation operator.

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1. concatenation of local basis matrices:

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and orthogonalization (truncation), using, e.g., SVD;

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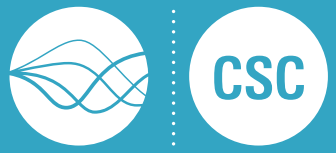
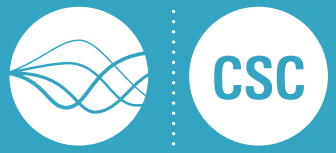
Avoids most of the problems encountered with local bases, but requires parameter-affine representation of system.



Using a Global Basis w/o Affine Parametrization

Empirical Matrix Interpolation Method [B./Gugercin/Willcox 2015]

Given $V, W \in \mathbb{R}^{n \times r}$ and suppose only that $M(p) \in \mathbb{R}^{n \times t}$ can be evaluated at specific parameter values.



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- Goal: approximate $m(p) \approx \tilde{m}(p) = \Psi\alpha(p)$, where $\Psi \in \mathbb{R}^{nt\times \ell}$ and $\alpha(p) \in \mathbb{R}^{\ell}$ with $\ell \ll n$.

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- Then $\hat{m}(p) = \operatorname{vec}(\hat{M}(p)) \in \mathbb{R}^{rt}$ (or $\mathbb{R}^{r^2}$ if $t = n$) can be computed cheaply and independent of n as

$$\begin{aligned}\hat{m}(p) &= \operatorname{vec}\left(W^T M(p) V\right)\\ &= (V^T \otimes W^T) m(p) \approx (V^T \otimes W^T)\tilde{m}(p) = (V^T \otimes W^T)\Psi\alpha(p) = \tilde{\tilde{m}}(p).\end{aligned}$$

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- Let $m(p) := \operatorname{vec}(M(p)) \in \mathbb{R}^{nt}$.
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- Then $\hat{m}(p) = \operatorname{vec}(\hat{M}(p)) \in \mathbb{R}^{rt}$ (or $\mathbb{R}^{r^2}$ if $t = n$) can be computed cheaply and independent of n as

$$\begin{aligned}\hat{m}(p) &= \operatorname{vec}\left(W^T M(p) V\right)\\ &= (V^T \otimes W^T) m(p) \approx (V^T \otimes W^T)\tilde{m}(p) = (V^T \otimes W^T)\Psi\alpha(p) = \tilde{\hat{m}}(p).\end{aligned}$$

- This is achieved by sampling $M(p)$ at $p = p^{(j)}$, $j = 1, \ldots, \ell$, yielding

$$\psi_j = \operatorname{vec}(M(p^{(j)})) \qquad \text{and} \qquad \Psi = [\psi_1, \ldots, \psi_\ell].$$

Then apply (Q,D)EIM (or alike) to determine $\alpha(p)$ s.t. selected entries of $\tilde{m}(p)$ interpolate those entries of $m(p)$.

Using a Global Basis w/o Affine Parametrization

Empirical Matrix Interpolation Method [B./Gugercin/Willcox 2015]

Given $V, W \in \mathbb{R}^{n\times r}$ and suppose only that $M(p) \in \mathbb{R}^{n\times t}$ can be evaluated at specific parameter values.

Goal: approximate $m(p) \approx \tilde{m}(p) = \Psi\alpha(p)$, where $\Psi \in \mathbb{R}^{nt\times \ell}$ and $\alpha(p) \in \mathbb{R}^{\ell}$ with $\ell \ll n$, and where Ψ is the sampling matrix built by $\operatorname{vec}(M(p^{(j)}))$.

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  Let $z_1, z_2, \ldots, z_\ell$ be the selected indices to be exactly matched, and $Z := [e_{z_1}, \ldots, e_{z_\ell}]$. Then, forcing interpolation at the selected rows implies

$$Z^T m(p) = Z^T \Psi\alpha(p) \quad \Longrightarrow \quad \alpha(p) = (Z^T\Psi)^{-1} Z^T m(p).$$

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$$\hat{M}(p) := \operatorname{vec}^{-1}\left(\tilde{\tilde{m}}(p)\right) = \operatorname{vec}^{-1}\left((V^T \otimes W^T)\Psi\alpha(p)\right) = \sum_{j=1}^{\ell} \alpha_j(p)\, \underbrace{W^T M(p^{(j)}) V}_{\text{precomputable!}}$$

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Error Measures

Parametric Systems Norms

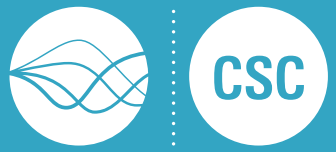
Mean-square error norm:

$$\|G - \hat{G}\|^2_{\mathcal{H}_2 \otimes L_2(\Omega)} := \frac{1}{2\pi} \int_{-\infty}^{+\infty} \int_{\Omega} \left\| G(\jmath\omega,\, p) - \hat{G}(\jmath\omega,\, p) \right\|_F^2 \, \mathrm{d}p_1 \ldots \mathrm{d}p_d \, \mathrm{d}\omega,$$

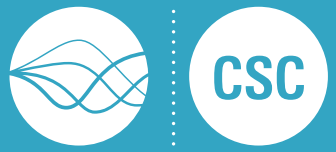
where $\|\,.\,\|_F$ denotes the Frobenius norm.

Worst-case error norm:

$$\|G - \hat{G}\|_{\mathcal{H}_\infty \otimes L_\infty(\Omega)} := \sup_{\omega \in \mathbb{R},\ p \in \Omega} \left\| G(\jmath\omega,\, p) - \hat{G}(\jmath\omega,\, p) \right\|_2 .$$



Outline



Recall Interpolatory Model Reduction

Computation of reduced-order model by projection

Given a linear (descriptor) system $E\dot{x} = Ax + Bu$, $y = Cx$ with transfer function $G(s) = C(sE - A)^{-1}B$, a reduced-order model is obtained using truncation matrices $V, W \in \mathbb{R}^{n\times r}$ with $W^TV = I_r$ ($\rightsquigarrow$ $(VW^T)^2 = VW^T$ is projector) by computing

$$\hat{E} = W^TEV,\ \hat{A} = W^TAV,\ \hat{B} = W^TB,\ \hat{C} = CV.$$

Petrov-Galerkin-type (two-sided) projection: $W \neq V$,

Galerkin-type (one-sided) projection: $W = V$.

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Rational Interpolation/Moment-Matching

Choose V, W such that

$$G(s_j) = \hat{G}(s_j), \quad j = 1, \ldots, k,$$

and

$$\frac{d^i}{ds^i} G(s_j) = \frac{d^i}{ds^i} \hat{G}(s_j), \quad i = 1, \ldots, K_j, \quad j = 1, \ldots, k.$$

Recall Interpolatory Model Reduction

Theorem (simplified) [Grimme 1997, Villemagne/Skelton 1987]

If

$$\begin{aligned} \operatorname{span}\left\{(s_1E - A)^{-1}B, \ldots, (s_kE - A)^{-1}B\right\} &\subset \operatorname{range}(V), \\ \operatorname{span}\left\{(s_1E - A)^{-T}C^T, \ldots, (s_kE - A)^{-T}C^T\right\} &\subset \operatorname{range}(W), \end{aligned}$$

then

$$G(s_j) = \hat{G}(s_j), \quad \frac{d}{ds}G(s_j) = \frac{d}{ds}\hat{G}(s_j), \quad \text{for } j = 1, \ldots, k.$$

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Remarks:

computation of V, W from rational Krylov subspaces, e.g.,

- dual rational Arnoldi/Lanczos [Grimme 1997],
- Iter. Rational Krylov-Alg. (IRKA) [Antoulas/Beattie/Gugercin 2006/08].

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Remarks:

using Galerkin/one-sided projection ($W \equiv V$) yields $G(s_j) = \hat{G}(s_j)$, but in general

$$\frac{d}{ds}G(s_j) \neq \frac{d}{ds}\hat{G}(s_j).$$

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Remarks:

$k = 1$, standard Krylov subspace(s) of dimension K:

$$\operatorname{range}(V) = \mathcal{K}_K((s_1E - A)^{-1}, (s_1E - A)^{-1}B).$$

$\rightsquigarrow$ **moment-matching methods**/Padé approximation,

$$\frac{d^i}{ds^i}G(s_1) = \frac{d^i}{ds^i}\hat{G}(s_1), \quad i = 0, \ldots, K-1(+K).$$

Comparison of Moment Matching and RBM

Numerical Example: A Printed Circuit Board (PCB)

- System in time domain:

$$\begin{aligned} E\dot{x}(t) &= Ax(t) + Bu(t), \\ y(t) &= Cx(t). \end{aligned}$$

- System in frequency domain:

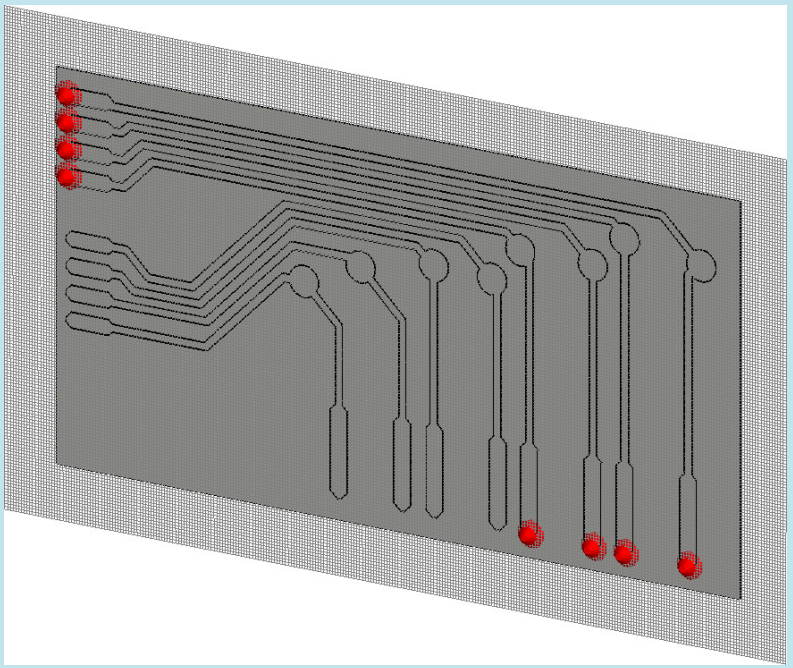
$$\begin{aligned} sEx(s) &= Ax(s) + Bu(s), \\ y(s) &= Cx(s). \end{aligned}$$

- Reduced basis method considers s as a parameter, and uses the system in frequency domain to compute

$$\mathrm{range}(V) = span\{x(s_1), \ldots, x(s_m)\}.$$

The ROM is obtained by Galerkin projection with V.

Printed circuit board



$n = 233,060,\ m = q = 1.$

Courtesy of TEMF, TU Darmstadt.

Comparison of Moment Matching and RBM

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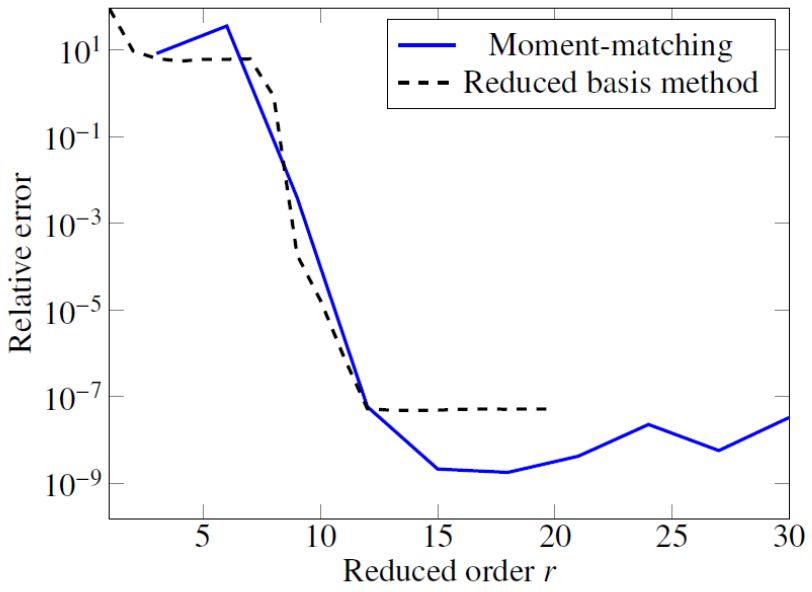
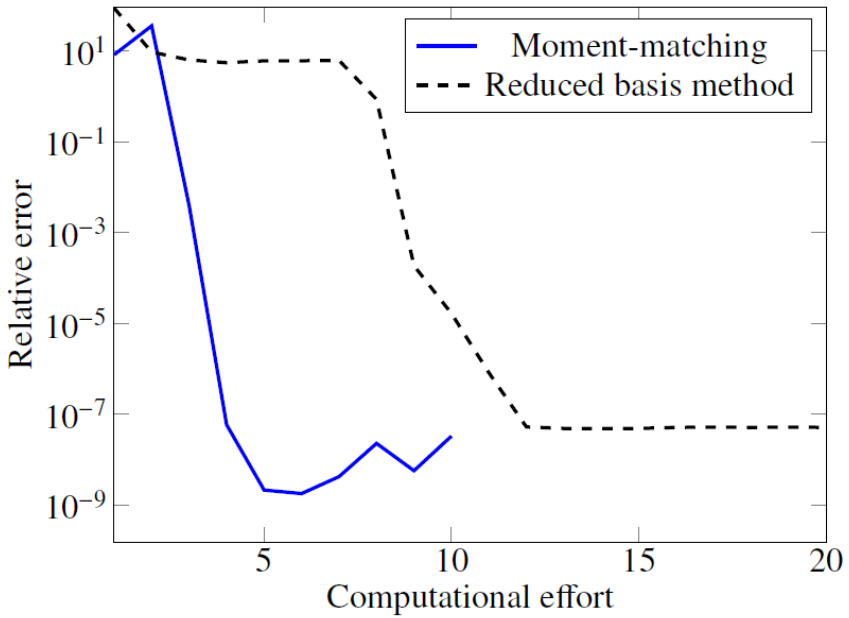
Moment-matching vs. reduced basis method





PMOR based on Multi-Moment Matching

Idea: choose appropriate frequency parameter $\hat{s}$ and parameter vector $\hat{p}$, expand into multivariate power series about $(\hat{s}, \hat{p})$ and compute reduced-order model, so that

$$G(s,p) = \hat{G}(s,p) + \mathcal{O}\left(|s-\hat{s}|^K + \|p-\hat{p}\|^L + |s-\hat{s}|^k \|p-\hat{p}\|^l\right),$$

i.e., first $K, L, k+l$ (mostly: $K = L = k+l$) coefficients (multi-moments) of Taylor/Laurent series coincide.

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Algorithms:

[1] [Daniel et al. 2004]: explicit computation of moments, numerically unstable.

[2] [Farle et al. 2006/07]: Krylov subspace approach, only polynomial param.-dependance, numerical properties not clear, but appears to be robust.

[3] [Weile et al. 1999, Feng/B. 2007/14]: Arnoldi-MGS method, employ recursive dependance of multi-moments, numerically robust, r often larger as for [2].

[4] **New:** employ dual-weighted residual error bound and greedy procedure to define interpolation points an # of multi-moments matched

[Antoulas/B./Feng 2014/17].

PMOR based on Multi-Moment Matching

Parametric System

Again, consider linear parametric system

$$\Sigma(p): \begin{cases} E(p)\dot{x}(t;p) &= A(p)x(t;p) + B(p)u(t), \\ y(t;p) &= C(p)x(t;p) \end{cases}$$

together with its transfer function $G(s,p)$.

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For simplicity, assume $B(\mu) \equiv B$, and re-parameterize — $\mu := [s, p^T, \dots]^T \in \mathbb{C}^\ell$ such that with

$$G(\mu) \equiv G(s,p), \quad x(\mu) \equiv x(s,p), \quad y(\mu) \equiv y(s,p), \dots$$
$$\mathcal{A}(\mu) := sE(p) - A(p),$$

we obtain linear-affine structure of $\mathcal{A}(\mu)$:

$$\mathcal{A}(\mu) = \mathcal{A}_0 + \mu_1 \mathcal{A}_1 + \dots + \mu_\ell \mathcal{A}_\ell.$$

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In frequency domain, we may then re-write the parametric system as

$$\mathcal{A}(\mu)x(\mu) = Bu(s), \qquad y(\mu) = C(\mu)x(\mu).$$

Multivariate Power Series Expansion I

Choose an expansion point $\mu^{(0)}$, and write

$$\begin{aligned}\mathcal{A}(\mu) &= \mathcal{A}_0 + \mu_1\mathcal{A}_1 + \ldots + \mu_\ell\mathcal{A}_m \\ &= \underbrace{(\mathcal{A}_0 + \mu_1^{(0)}\mathcal{A}_1 + \ldots + \mu_\ell^{(0)}\mathcal{A}_m)}_{:=\mathcal{M}_0} + \Big((\mu_1 - \mu_1^{(0)})\mathcal{A}_1 + \ldots + (\mu_\ell - \mu_\ell^{(0)})\mathcal{A}_\ell\Big)\end{aligned}$$

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Using the Neumann lemma ($(I-F)^{-1} = \sum_{j=0}^{\infty} F^j$ if $\|F\| < 1$), we obtain

$$\begin{aligned}\mathcal{A}(\mu)^{-1} &= \sum_{j=0}^{\infty}(-1)^j\left((\mu_1 - \mu_1^{(0)})\mathcal{M}_0^{-1}\mathcal{A}_1 + \ldots + (\mu_\ell - \mu_\ell^{(0)})\mathcal{M}_0^{-1}\mathcal{A}_\ell\right)^j\mathcal{M}_0^{-1} \\ &= \sum_{j=0}^{\infty}(\sigma_1\mathcal{M}_1 + \ldots\sigma_\ell\mathcal{M}_\ell)^j\mathcal{M}_0^{-1},\end{aligned}$$

where $\sigma_i = \mu_i - \mu_i^{(0)}$ and $\mathcal{M}_i = -\mathcal{M}_0^{-1}\mathcal{A}_i$ for $i = 1, \ldots, \ell$.

Multivariate Power Series Expansion II

We have

$$\mathcal{A}(\mu)x(\mu) = Bu(s).$$

and

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Hence,

$$\begin{aligned} x(\mu) = \mathcal{A}(\mu)^{-1}Bu(s) &= \sum_{j=0}^{\infty}(\sigma_1\mathcal{M}_1 + \ldots + \sigma_\ell\mathcal{M}_\ell)^j \underbrace{\mathcal{M}_0^{-1}B}_{=:\mathcal{B}} u(s) \\ &\approx \sum_{j=0}^{k}(\sigma_1\mathcal{M}_1 + \ldots + \sigma_\ell\mathcal{M}_\ell)^j\mathcal{B}u(s) =: \tilde{x}(\mu). \end{aligned}$$

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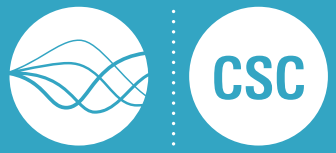
Thus, $x(\mu)$ is (approximately, locally) contained in the Krylov subspace $\mathcal{K}_{k+1}((\sigma_1\mathcal{M}_1 + \ldots + \sigma_\ell\mathcal{M}_\ell), \mathcal{B})$.



PMOR based on Multi-Moment Matching

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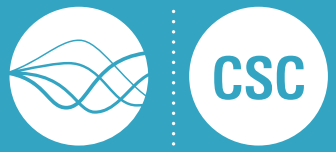
- Project the state-space onto this subspace.



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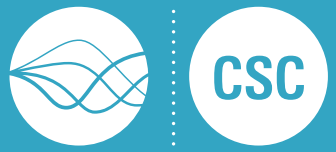

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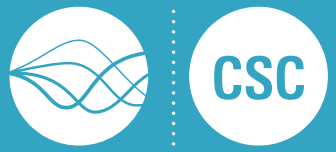

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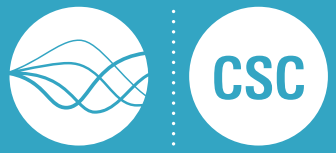

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- First terms in the multivarite Taylor expansion match, i.e., we achieve matrix interpolation for partial derivatives up to order ℓ, or more in the Petrov-Galerkin case.
- Approximation is only valid locally (convergence radius of Neumann series!) $\rightsquigarrow$ use several expansion points $\mu^{(0)}, \ldots, \mu^{(h)}$, and concatenate (and truncate) the local bases to obtain a global basis.

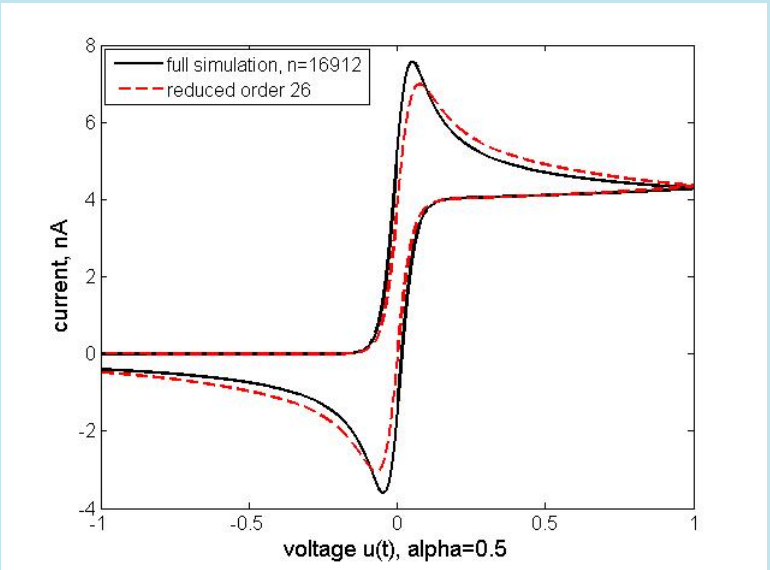
Numerical Examples: Electro-Chemical SEM

Compute cyclic voltammogram based on FE model

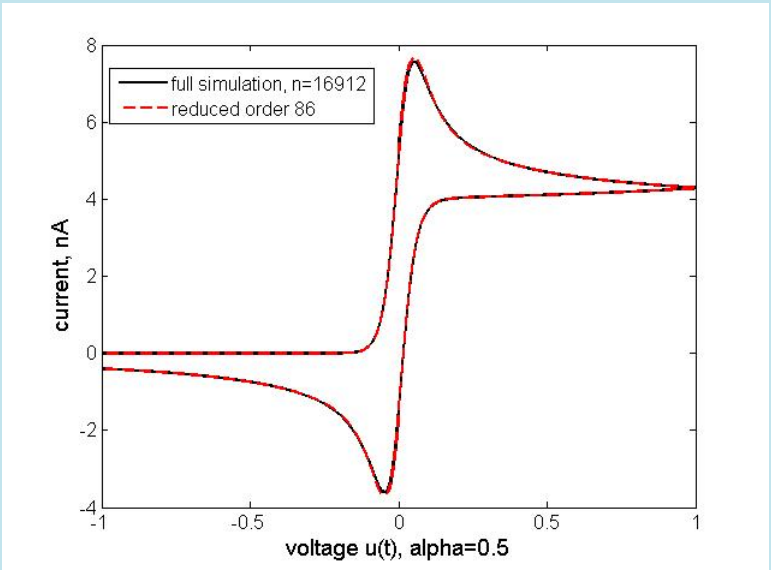
$$E\dot{x}(t) = (A_0 + p_1 A_1 + p_2 A_2)x(t) + Bu(t), \quad y(t) = c^T x(t),$$

where $n = 16,912$, $m = 3$, A_1, A_2 diagonal.

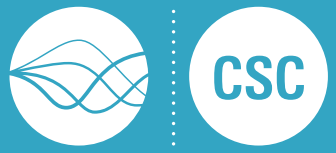
$K = L = k + \ell = 4 \;\Rightarrow\; r = 26$



$K = L = k + \ell = 9 \;\Rightarrow\; r = 86$



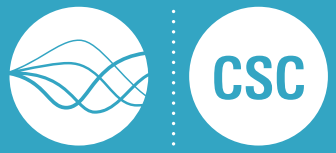
Source: MOR Wiki: http://morwiki.mpi-magdeburg.mpg.de/morwiki/index.php/Scanning_Electrochemical_Microscopy



PMOR based on Multi-Moment Matching

Open question

How to adaptively choose $\mu^{(i)}$?



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And how many partial derivatives to be matched at each interpolation point?

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And how many partial derivatives to be matched at each interpolation point?

Possible approach: adopt ideas from Reduced Basis Methods, i.e., let

$$\|G(\mu) - \hat{G}(\mu)\| \leq \Delta(\mu) \quad \text{or} \quad \|y(\mu) - \hat{y}(\mu)\| \leq \Delta_o(\mu)$$

guide the selection of $\mu^{(i)}$ for computable *a posteriori* error bounds for the state or the output.

Error Bound for Automatic ROM Construction

Theorem (SISO case) [Feng/Antoulas/B. 2015/17]

Assume that $\sigma_{\min}(G(s,p)) =: \beta(s,p) > 0 \quad \forall\, \mathrm{Re}(s) \geq 0, \forall\, p \in \Omega$, then

$$|H(s,p) - \hat{H}(s,p)| \leq \tilde{\Delta}(s,p) + |(\hat{x}^{du})^H \mathbf{r}^{pr}(s,p)| =: \Delta(s,p),$$

where

$$\tilde{\Delta}(s,p) = \frac{\|r^{du}(s,p)\|_2 \|r^{pr}(s,p)\|_2}{\beta(s,p)},$$

with the primal and dual residuals r^{pr}, r^{du} and the reduced "dual state" $\hat{x}^{du}$:

$$\begin{aligned} r^{pr}(s,p) &= \| \left(B - (sE(p) - A(p))\right) \left(V(s\hat{E}(p) - \hat{A}(p))^{-1} \hat{B}\right) \|, \\ r^{du}(s,p) &= \| \left(C^T - (\bar{s}E(p) - A(p))^T\right) \hat{x}^{du} \|, \\ \hat{x}^{du} &= -V^{du}(\bar{s}\hat{E}^{du}(p) - \hat{A}^{du}(p))^{-T} \hat{C}^{du}. \end{aligned}$$

The dual reduced-order system is computed using Galerkin projection with V^{du} obtained by applying multi-moment matching algorithm to "dual" system $(\bar{s}E(p)^T - A(p)^T, C^T)$.

Error Bound for Automatic ROM Construction

Remarks

- For application in "RBM fashion", $r^{du}(\mu), r^{pr}(\mu)$ can be efficiently computed, need to solve sparse linear systems on training set, i.e., one sparse factorization for each sampling point.
- $\beta(s,p) = \sigma_{\min}(G(s,p))$ easily computable on the training set — system solves for evaluation of the transfer function readily available from residual computation!
- Extension to MIMO case possible taking max over all I/O channels.
- Can use Petrov-Galerkin framework using $W = V^{du}$ at no extra cost!

Algorithm 1 Automatic generation of the ROM: adaptively selecting $\mu^{(i)}$

Input: $V = [\,]$; $\epsilon > \epsilon_{tol}$; Initial expansion point: $\hat{\mu}$; $i := -1$;
Ξ_{train}: a set of samples of μ covering the parameter domain.

Output: V.

1: **while** $\epsilon > \epsilon_{tol}$ **do**
2: $\quad i = i + 1$;
3: $\quad \mu^{(i)} = \hat{\mu}$;
4: $\quad V_{\mu^{(i)}} =$ orthogonal basis of $\mathcal{K}_{k+1}((\sigma_1^{(i)}\mathcal{M}_1 + \ldots + \sigma_\ell^{(i)}\mathcal{M}_\ell), \mathcal{B})$;
5: $\quad V = \mathtt{orth}([V, V_{\mu^{(i)}}])$;
6: $\quad \hat{\mu} = \arg \max_{\mu \in \Xi_{train}} \Delta(\mu)$;
7: $\quad \epsilon = \Delta(\hat{\mu})$;
8: **end while**

Numerical Example: Silicon Nitride Membrane

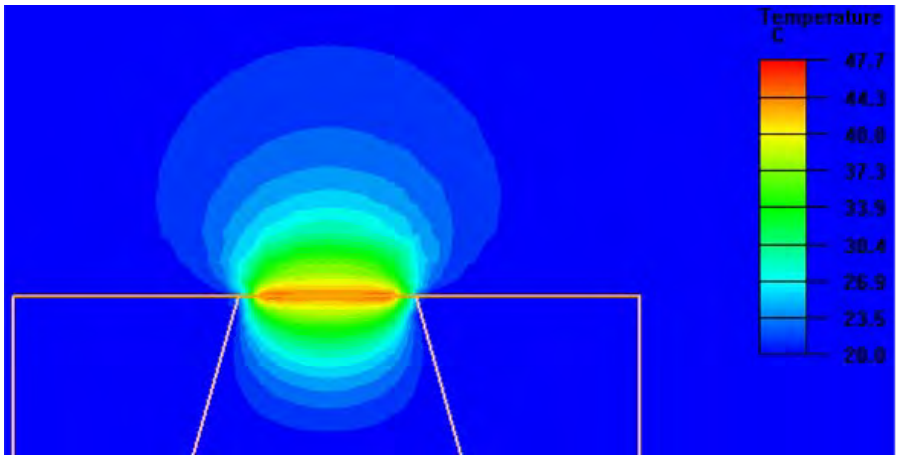
A SiN membrane can be a part of a gas sensor, an infra-red sensor, a microthruster, etc. Heat tansfer in the membrane is described by

$$\begin{aligned}(E_0 + \rho c_p E_1)\dot{x}(t) &= -(K_0 + \kappa K_1 + hK_2)x(t) + bu(t)\\ y(t) &= Cx(t),\end{aligned}$$

with parameters

- density $\rho \in [3000, 3200]$,
- specific heat capacity $c_p \in [400, 750]$,
- thermal conductivity $\kappa \in [2.5, 4]$,
- membrane heat transfer coefficient $h \in [10, 12]$.

and frequency $f \in [0, 25]Hz$.



Source: MOR Wiki: http://morwiki.mpi-magdeburg.mpg.de/morwiki/index.php/Silicon_nitride_membrane

Numerical Examples: Silicon Nitride Membrane

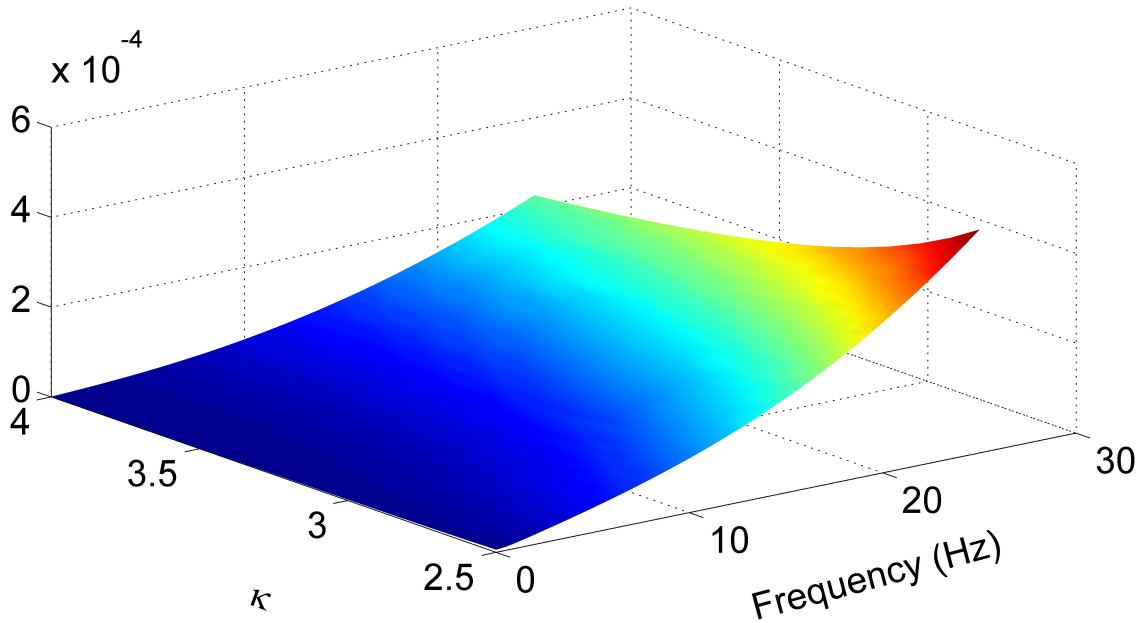
Setting

- Training set: $\Xi_{train} = 5$ random samples for ρ and c_p, 3 random samples for κ and h, respectively, 10 samples of Laplace variable s.
- Error measures:
  $\varepsilon^{rel}_{true} = \max\limits_{\mu\in\Xi_{train}} |G(\mu) - \hat{G}(\mu)|/|G(\mu)|,$
  $\Delta^{rel}(\mu) = \Delta(\mu)/|\hat{G}(\mu)|$
- $V_{\mu^{(i)}} = \mathrm{span}\{\mathcal{B}, (\sigma_1^{(i)}\mathcal{M}_1 + \ldots + \sigma_\ell^{(i)}\mathcal{M}_\ell)\mathcal{B}\}$, $\epsilon^{re}_{tol} = 10^{-2}$, $n = 60,020$, $r = 8$.

iter.	ε^{rel}_{true}	$\Delta^{rel}(\mu^{(i)})$	s	ρc_p	κ	h
1	1×10^{-3}	3.44	18.94	1.37×10^{6}	2.74	10.97
2	1×10^{-4}	4.59×10^{-2}	0.89	1.31×10^{6}	3.96	11.60
3	2.80×10^{-5}	4.07×10^{-2}	23.98	2.35×10^{6}	3.94	10.28
4	2.58×10^{-6}	2.62×10^{-5}	0.89	2.31×10^{6}	2.74	10.28

Numerical Examples: Silicon Nitride Membrane

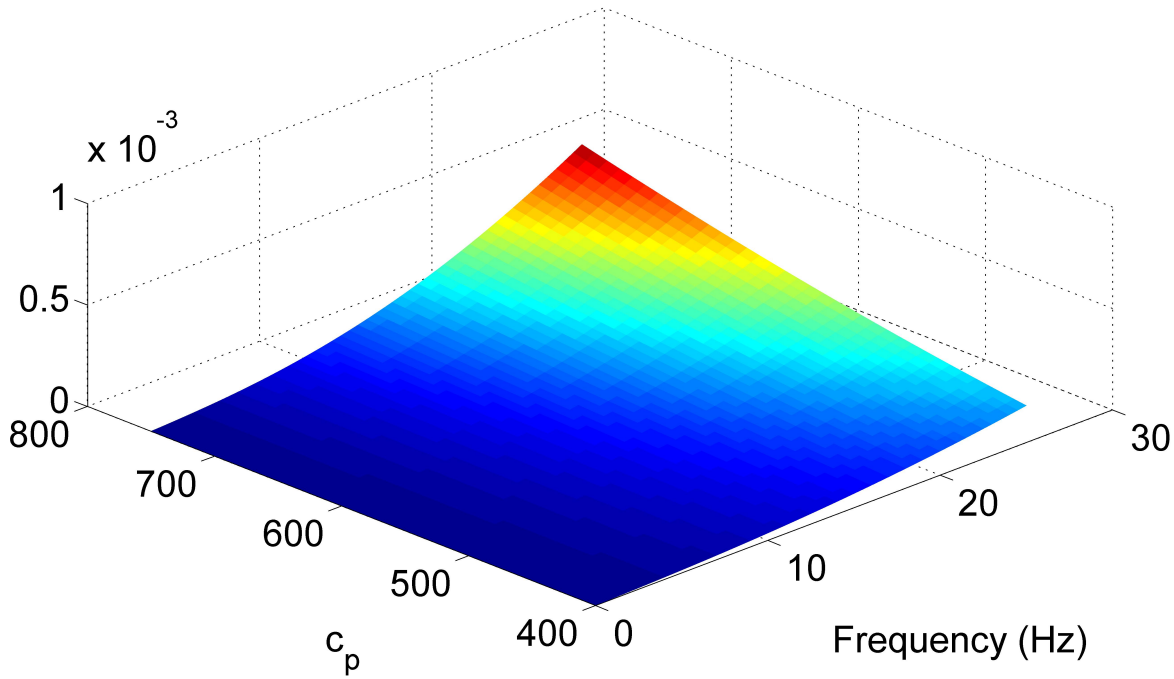
Verification of the accuracy of the ROM for κ over set Ξ_{fine} with 16 equidistant samples of κ, 51 equidistant samples of the frequency f, while the other parameters are fixed.



Relative error of the final ROM changing with κ and frequency.

Numerical Examples: Silicon Nitride Membrane

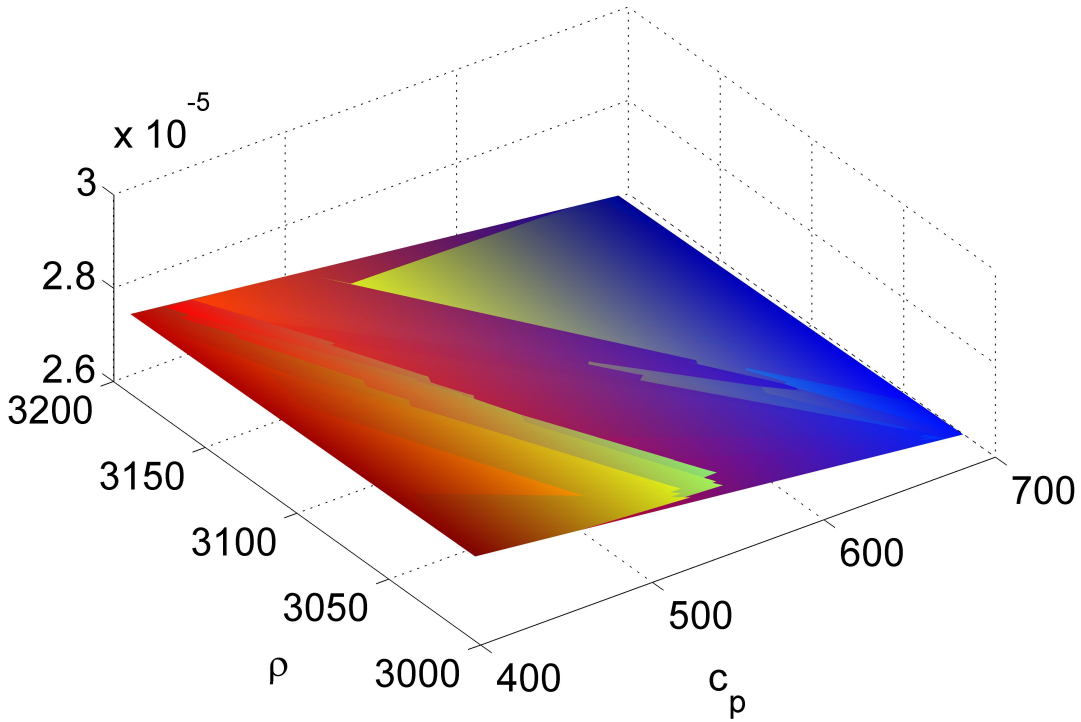
Verification of the accuracy of the ROM for c_p over set Ξ_{fine} with 36 equidistant samples of c_p, 51 equidistant samples of the frequency f, while the other parameters are fixed.



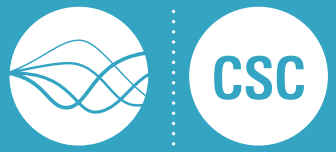
Relative error of the final ROM changing with c_p and frequency.

Numerical Examples: Silicon Nitride Membrane

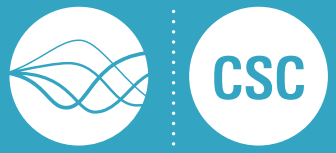
Verification of the accuracy of the ROM for ρ, c_p over set Ξ_{fine} with 50 random samples of ρ, c_p, respectively, the other parameters are fixed.



Relative error of the final ROM changing with c_p and κ.

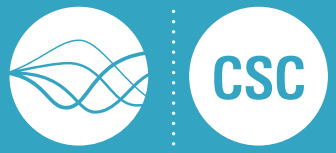


Outline



PMOR based on Multi-Moment Matching

Greedy expansion point selection has a heuristic nature and relies on a training set.



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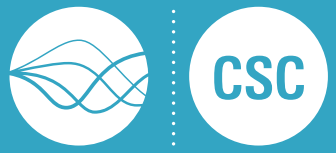

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- Iterative procedure for selecting interpolation points. [χ] ... [✓] for special case.

$\mathcal{H}_2$-Model Reduction for Linear Systems

Consider stable (i.e. $\Lambda(A) \subset \mathbb{C}^-$) linear systems Σ,

$$\dot{x}(t) = Ax(t) + Bu(t),\ y(t) = Cx(t) \quad \simeq \quad y(s) = \underbrace{C(sI - A)^{-1}B}_{=:G(s)}\, u(s)$$

System norms

Recall: two common system norms for measuring approximation quality are

- $\mathcal{H}_2$-norm, $\|\Sigma\|_{\mathcal{H}_2} = \left(\frac{1}{2\pi}\int_0^{2\pi} \operatorname{tr}\left(\left(G^T(-\jmath\omega)G(\jmath\omega)\right)\right) d\omega\right)^{\frac{1}{2}}$,
- $\mathcal{H}_\infty$-norm, $\|\Sigma\|_{\mathcal{H}_\infty} = \sup\limits_{\omega\in\mathbb{R}} \sigma_{\max}(G(\jmath\omega))$,

where

$$G(s) = C(sI - A)^{-1}B.$$

Error system and $\mathcal{H}_2$-Optimality [Meier/Luenberger 1967]

In order to find an $\mathcal{H}_2$-optimal reduced system, consider the error system $G(s) - \hat{G}(s)$ which can be realized by

$$A^{err} = \begin{bmatrix} A & 0 \\ 0 & \hat{A} \end{bmatrix}, \quad B^{err} = \begin{bmatrix} B \\ \hat{B} \end{bmatrix}, \quad C^{err} = \begin{bmatrix} C & -\hat{C} \end{bmatrix}.$$

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Assuming a coordinate system in which $\hat{A}$ is diagonal and taking derivatives of

$$\|G(\,.\,) - \hat{G}(\,.\,)\|_{\mathcal{H}_2}^2$$

with respect to free parameters in $\Lambda(\hat{A}), \hat{B}, \hat{C} \rightsquigarrow$ first-order necessary $\mathcal{H}_2$-optimality conditions (SISO)

$$G(-\hat{\lambda}_i) = \hat{G}(-\hat{\lambda}_i),$$
$$G'(-\hat{\lambda}_i) = \hat{G}'(-\hat{\lambda}_i),$$

where $\hat{\lambda}_i$ are the poles of the reduced system $\hat{\Sigma}$.

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First-order necessary $\mathcal{H}_2$-optimality conditions (MIMO):

$$\begin{aligned} G(-\hat{\lambda}_i)\tilde{B}_i &= \hat{G}(-\hat{\lambda}_i)\tilde{B}_i, && \text{for } i = 1, \ldots, r, \\ \tilde{C}_i^T G(-\hat{\lambda}_i) &= \tilde{C}_i^T \hat{G}(-\hat{\lambda}_i), && \text{for } i = 1, \ldots, r, \\ \tilde{C}_i^T H'(-\hat{\lambda}_i)\tilde{B}_i &= \tilde{C}_i^T \hat{G}'(-\hat{\lambda}_i)\tilde{B}_i && \text{for } i = 1, \ldots, r, \end{aligned}$$

where $\hat{A} = R\hat{\Lambda}R^{-T}$ is the spectral decomposition of the reduced system and $\tilde{B} = \hat{B}^T R^{-T}$, $\tilde{C} = \hat{C}R$.

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$$\begin{aligned} \Leftrightarrow \quad & \operatorname{vec}(I_q)^T \left(e_j e_i^T \otimes C\right) \left(-\hat{\Lambda} \otimes I_n - I_r \otimes A\right)^{-1} \left(\tilde{B}^T \otimes B\right) \operatorname{vec}(I_m) \\ & = \operatorname{vec}(I_q)^T \left(e_j e_i^T \otimes \hat{C}\right) \left(-\hat{\Lambda} \otimes I_r - I_r \otimes \hat{A}\right)^{-1} \left(\tilde{B}^T \otimes \hat{B}\right) \operatorname{vec}(I_m), \\ & \qquad \text{for } i = 1, \ldots, r \text{ and } j = 1, \ldots, q. \end{aligned}$$

Interpolation of the Transfer Function [Grimme 1997]

Construct reduced transfer function by Petrov-Galerkin projection $\mathcal{P} = VW^T$, i.e.

$$\hat{G}(s) = CV\left(sI - W^TAV\right)^{-1} W^TB,$$

Interpolation of the Transfer Function [Grimme 1997]

Construct reduced transfer function by Petrov-Galerkin projection $\mathcal{P} = VW^T$, i.e.

$$\hat{G}(s) = CV\left(sI - W^TAV\right)^{-1}W^TB,$$

where V and W are given as

$$V = \left[(-\mu_1 I - A)^{-1}B, \ldots, (-\mu_r I - A)^{-1}B\right],$$
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Then

$$G(-\mu_i) = \hat{G}(-\mu_i) \quad \text{and} \quad G'(-\mu_i) = \hat{G}'(-\mu_i),$$

for $i = 1, \ldots, r$.

$\mathcal{H}_2$-optimal Model Reduction for Linear Systems

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Then

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for $i = 1, \ldots, r$.
Starting with an initial guess for $\hat{\Lambda}$ and setting $\mu_i \equiv \hat{\lambda}_i \rightsquigarrow$ iterative algorithms (IRKA/MIRIAm) that yield $\mathcal{H}_2$-optimal models.

[Gugercin et al. 2006/08], [Bunse-Gerstner et al. 2007], [Van Dooren et al. 2008]

The Basic IRKA Algorithm

Algorithm 2 IRKA (MIMO version/MIRIAm)

Input: A stable, B, C, $\hat{A}$ stable, $\hat{B}$, $\hat{C}$, $\delta > 0$.
Output: A^{opt}, B^{opt}, C^{opt}

1: **while** $\left(\max_{j=1,\ldots,r} \left\{ \frac{|\mu_j - \mu_j^{\text{old}}|}{|\mu_j|} \right\} > \delta\right)$ **do**
2: $\quad \operatorname{diag}(\mu_1, \ldots, \mu_r) := R^{-1}\hat{A}R =$ spectral decomposition.
3: $\quad \tilde{B} = \hat{B}^H R^{-T}$, $\tilde{C} = \hat{C}R$.
4: $\quad V = \left[(-\mu_1 I - A)^{-1} B\tilde{b}_1, \ldots, (-\mu_r I - A)^{-1} B\tilde{b}_r\right]$
5: $\quad W = \left[(-\mu_1 I - A^T)^{-1} C^T\tilde{c}_1, \ldots, (-\mu_r I - A^T)^{-1} C^T\tilde{c}_r\right]$
6: $\quad V = \operatorname{orth}(V)$, $W = \operatorname{orth}(W)$, $W = W(V^H W)^{-1}$
7: $\quad \hat{A} = W^H AV$, $\hat{B} = W^H B$, $\hat{C} = CV$.
8: **end while**
9: $A^{opt} = \hat{A}$, $B^{opt} = \hat{B}$, $C^{opt} = \hat{C}$.

Theory: Interpolation of the Transfer Function

Theorem [Baur/Beattie/B./Gugercin 2007/11]

Let
$$\begin{aligned}\hat{G}(s,p) &:= \hat{C}(p)(s\hat{E}(p)-\hat{A}(p))^{-1}\hat{B}(p)\\ &= C(p)V(sW^TE(p)V-W^TA(p)V)^{-1}W^TB(p).\end{aligned}$$

Suppose $\hat{p}=[\hat{p}_1,\ldots,\hat{p}_d]^T$ and $\hat{s}\in\mathbb{C}$ are chosen such that both $\hat{s}\,E(\hat{p})-A(\hat{p})$ and $\hat{s}\,\hat{E}(\hat{p})-\hat{A}(\hat{p})$ are invertible. If

$$(\hat{s}\,E(\hat{p})-A(\hat{p}))^{-1}\,B(\hat{p})\in\operatorname{range}(V)$$

or

$$\left(C(\hat{p})\,(\hat{s}\,E(\hat{p})-A(\hat{p}))^{-1}\right)^T\in\operatorname{range}(W),$$

then $G(\hat{s},\hat{p})=\hat{G}(\hat{s},\hat{p})$.

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then $G(\hat{s},\hat{p})=\hat{G}(\hat{s},\hat{p})$.

Extension to MIMO case using tangential interpolation: let $0\neq b\in\mathbb{R}^m$, $0\neq c\in\mathbb{R}^q$.

a) If $(\hat{s}\,E(\hat{p})-A(\hat{p}))^{-1}\,B(\hat{p})b\in\operatorname{range}(V)$, then $G(\hat{s},\hat{p})b=\hat{G}(\hat{s},\hat{p})b$.

b) If $\left(c^TC(\hat{p})\,(\hat{s}\,E(\hat{p})-A(\hat{p}))^{-1}\right)^T\in\operatorname{range}(W)$, then $c^TG(\hat{s},\hat{p})=c^T\hat{G}(\hat{s},\hat{p})$.

Theory: Interpolation of the Parameter Gradient

Theorem [Baur/Beattie/B./Gugercin 2007/11]

Suppose that $E(p)$, $A(p)$, $B(p)$, $C(p)$ are C^1 in a neighborhood of $\hat{p} = [\hat{p}_1, \ldots, \hat{p}_d]^T$ and that both $\hat{s}\,E(\hat{p}) - A(\hat{p})$ and $\hat{s}\,\hat{E}(\hat{p}) - \hat{A}(\hat{p})$ are invertible.

If

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then

$$\nabla_p G(\hat{s}, \hat{p}) = \nabla_p G_r(\hat{s}, \hat{p}), \qquad \frac{\partial}{\partial s} G(\hat{s}, \hat{p}) = \frac{\partial}{\partial s} \hat{G}(\hat{s}, \hat{p}).$$

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Note: result extends to MIMO case using tangential interpolation:

Let $0 \neq b \in \mathbb{R}^m$, $0 \neq c \in \mathbb{R}^q$ be arbitrary. If $(\hat{s}\,E(\hat{p}) - A(\hat{p}))^{-1} B(\hat{p}) b \in \operatorname{range}(V)$ and $\left(c^T C(\hat{p})\,(\hat{s}\,E(\hat{p}) - A(\hat{p}))^{-1}\right)^T \in \operatorname{range}(W)$, then $\nabla_p c^T G(\hat{s}, \hat{p}) b = \nabla_p c^T \hat{G}(\hat{s}, \hat{p}) b$.

$\mathcal{H}_2$-(sub)optimal Model Reduction for Linear Parametric Systems

Theory: Interpolation of the Parameter Gradient

Theorem [Baur/Beattie/B./Gugercin 2007/11]

Suppose that $E(p)$, $A(p)$, $B(p)$, $C(p)$ are C^1 in a neighborhood of $\hat{p} = [\hat{p}_1, \ldots, \hat{p}_d]^T$ and that both $\hat{s}\,E(\hat{p}) - A(\hat{p})$ and $\hat{s}\,\hat{E}(\hat{p}) - \hat{A}(\hat{p})$ are invertible.

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1. Reduced-order model satisfies necessary conditions for surrogate models in trust region methods [Alexandrov/Dennis/Lewis/Torczon 1998].
2. Approximation of gradient allows use of reduced-order model for sensitivity analysis.

Generic implementation of interpolatory PMOR

Define $\mathcal{A}(s,p) := sE(p) - A(p)$.

1. Select "frequencies" $s_1, \ldots, s_k \in \mathbb{C}$ and parameter vectors $p^{(1)}, \ldots, p^{(\ell)} \in \Omega \subset \mathbb{R}^d$.
2. Compute (orthonormal) basis of
$$\mathcal{V} = \operatorname{span}\left\{\mathcal{A}(s_1, p^{(1)})^{-1} B(p^{(1)}), \ldots, \mathcal{A}(s_k, p^{(\ell)})^{-1} B(p^{(\ell)})\right\}.$$
3. Compute (orthonormal) basis of
$$\mathcal{W} = \operatorname{span}\left\{\mathcal{A}(s_1, p^{(1)})^{-T} C(p^{(1)})^T, \ldots, \mathcal{A}(s_k, p^{(\ell)})^{-T} C(p^{(\ell)})^T\right\}.$$
4. Set $V := [v_1, \ldots, v_{k\ell}]$, $\tilde{W} := [w_1, \ldots, w_{k\ell}]$, and $W := \tilde{W}(\tilde{W}^T V)^{-1}$. (Note: $r = k\ell$).
5. Compute $\begin{cases} \hat{A}(p) := W^T A(p) V, & \hat{B}(p) := W^T B(p) V, \\ \hat{C}(p) := W^T C(p) V, & \hat{E}(p) := W^T E(p) V. \end{cases}$

Remarks

- If directional derivatives w.r.t. p are included in $\operatorname{range}(V)$, $\operatorname{range}(W)$, then also the Hessian of $G(\hat{s}, \hat{p})$ is interpolated by the Hessian of $\hat{G}(\hat{s}, \hat{p})$.

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- For prescribed parameter vectors $p^{(k)}$, we can use corresponding $\mathcal{H}_2$-optimal frequencies $s_{k,\ell}$, $\ell = 1,\ldots,r_k$ computed by IRKA, i.e., reduced-order systems $\hat{G}_*^{(k)}$ so that

$$\|G(.,p^{(k)}) - \hat{G}_*^{(k)}(.)\|_{\mathcal{H}_2} = \min_{\substack{\operatorname{order}(\hat{G})=r_k \\ \hat{G}\ \text{stable}}} \|G(.,p^{(k)}) - \hat{G}^{(k)}(.)\|_{\mathcal{H}_2},$$

where

$$\|G\|_{\mathcal{H}_2} := \left(\frac{1}{2\pi}\int_{-\infty}^{+\infty} \|G(\jmath\omega)\|_{\mathrm{F}}^2 d\omega\right)^{1/2}.$$

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- Optimal choice of interpolation frequencies s_k and parameter vectors $p^{(k)}$ possible for special cases.

Numerical Example: Thermal Conduction in a Semiconductor Chip

- Important requirement for a compact model of thermal conduction is boundary condition independence.
- The thermal problem is modeled by the heat equation, where heat exchange through device interfaces is modeled by convection boundary conditions containing film coefficients $\{p_i\}_{i=1}^3$ describing the heat exchange at ith interface.
- Spatial semi-discretization leads to

$$E\dot{x}(t) = (A_0 + \sum_{i=1}^{3} p_i A_i)x(t) + bu(t), \quad y(t) = c^T x(t),$$

where $n = 4,257$, A_i, $i = 1, 2, 3$, are diagonal.

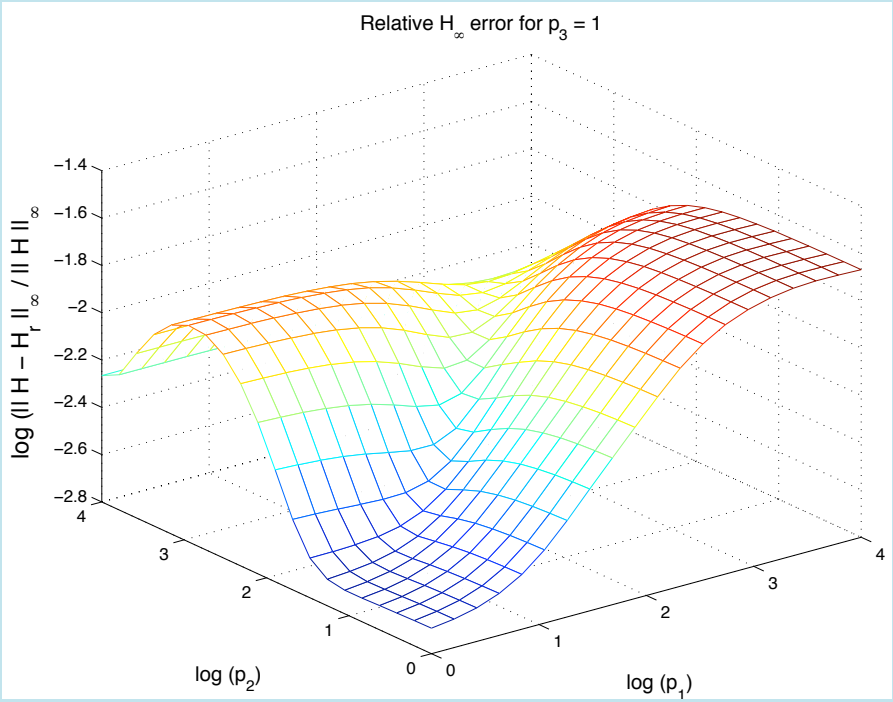
Source: C.J.M Lasance, *Two benchmarks to facilitate the study of compact thermal modeling phenomena*, IEEE Transactions on Components and Packaging Technologies, 24(4):559–565, 2001.

MOR Wiki: http://morwiki.mpi-magdeburg.mpg.de/morwiki/index.php/Microthruster_Unit

Numerical Example: Thermal Conduction in a Semiconductor Chip

Choose 2 interpolation points for parameters ("important" configurations), 8/7 H_2-optimal interpolation frequencies selected by IRKA. $\Rightarrow$ $k = 2, \ell = 8, 7$, hence $r = 15$.

$p_3 = 1$, $p_1, p_2 \in [1, 10^4]$.



Optimality of Interpolation Points

Theorem [Baur/Beattie/B./Gugercin 2011]

For special parameterized SISO systems,

$$A(p) \equiv A_0, \quad E(p) \equiv E_0, \quad B(p) = B_0 + p_1 B_1, \quad C(p) = C_0 + p_2 C_1,$$

optimal choice possible, necessary conditions:
If $\hat{G}$ minimizes the approximation error w.r.t. $\|G - \hat{G}\|_{\mathcal{H}_2 \times L_2(\Omega)}$, $p \in \Omega \subset \mathbb{R}^d$, and $\Lambda(\hat{A}, \hat{E}) = \{\hat{\lambda}_1, \ldots, \hat{\lambda}_r\}$ (all simple), then the interpolation frequencies satisfy

$$s_i = -\hat{\lambda}_i, \quad i = 1, \ldots, r,$$

and the parameter interpolation points $\{p^{(1)}, \ldots, p^{(r)}\}$ satisfy the interpolation conditions

$$\begin{aligned} G(-\hat{\lambda}_k, p^{(k)}) &= \hat{G}(-\hat{\lambda}, p^{(k)}), \\ \frac{\partial}{\partial s} G(-\hat{\lambda}, p^{(k)}) &= \frac{\partial}{\partial s} \hat{G}(-\hat{\lambda}, p^{(k)}), \quad \nabla_p G(-\hat{\lambda}, p^{(k)}) = \nabla_p \hat{G}(-\hat{\lambda}, p^{(k)}). \end{aligned}$$

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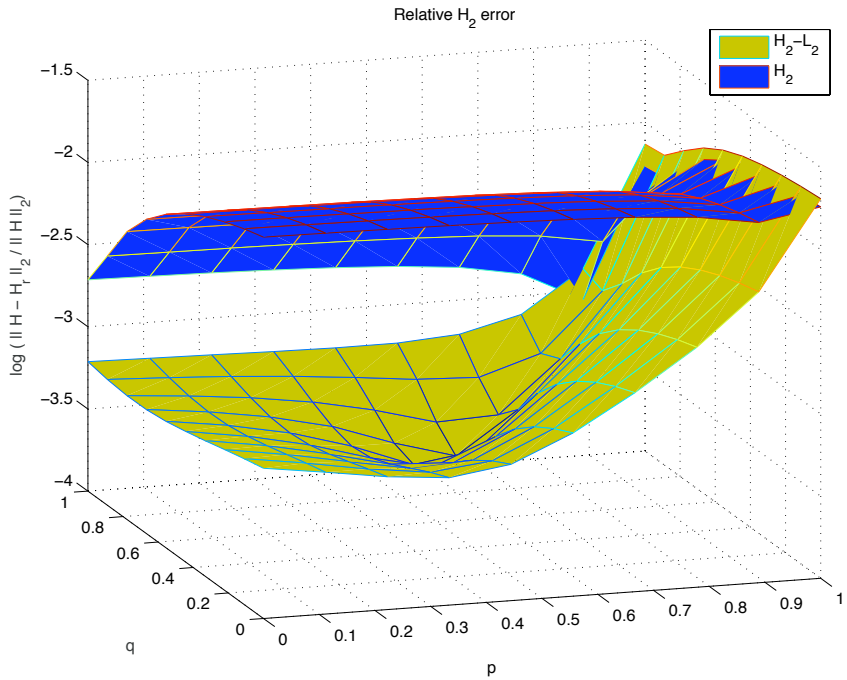
Proof:

$$\|G\|_{\mathcal{H}_2 \times L_2(\Omega)} = \|L^T \tilde{G} L\|_{\mathcal{H}_2}, \quad \text{where } \tilde{G}(s) = \begin{bmatrix} C_0 \\ C_1 \end{bmatrix} (sE - A)^{-1} [\, B_0, B_1 \,],\ L = \begin{bmatrix} 1 & 0 \\ \frac{1}{2} & \frac{1}{2\sqrt{3}} \end{bmatrix}.$$

$\Longrightarrow$ **Computation via IRKA applied to $\tilde{G}$.**

Optimality of Interpolation Points — Numerical Example

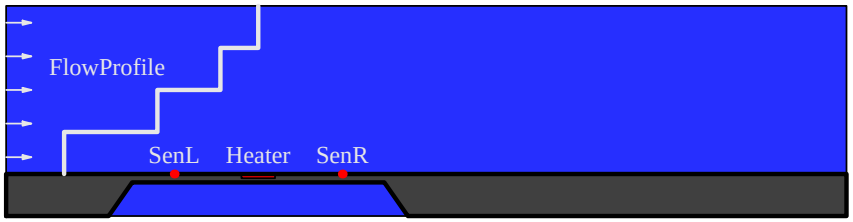
- Model for evolution of temperature distribution on a plate, described by the heat equation.
- FDM SISO model of order $n = 197$.
- Parameter $p_1 \in [0, 1]$ encodes movement of heat source from B_0 to $B_0 + B_1$, analogous for relocation of measurement.

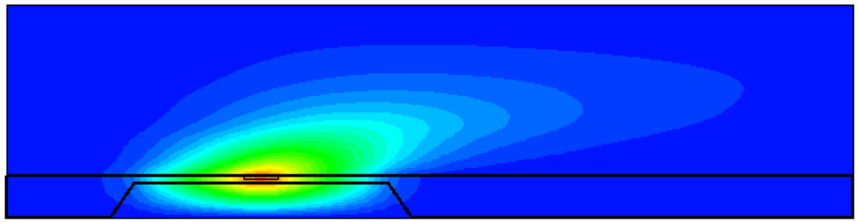


Relative $\mathcal{H}_2 \otimes L_2(\Omega)$ error: 7.5×10^{-4}.

A Comparison of PMOR Methods: Anemometer

Consider an anemometer, a flow sensing device located on a membrane used in the context of minimizing heat dissipation.





Source: [Baur/B./Greiner/Korvink/Lienemann/Moosmann 2011]

- FE model:

$$E\dot{x}(t) = (A + pA_1)x(t) + Bu(t), \quad y(t) = Cx(t), \quad x(0) = 0,$$

- $n = 29,008,\ m = 1,\ q = 3,\ p_1 \in [0, 1]$ fluid velocity.

Source: MOR Wiki: http://morwiki.mpi-magdeburg.mpg.de/morwiki/index.php/Anemometer

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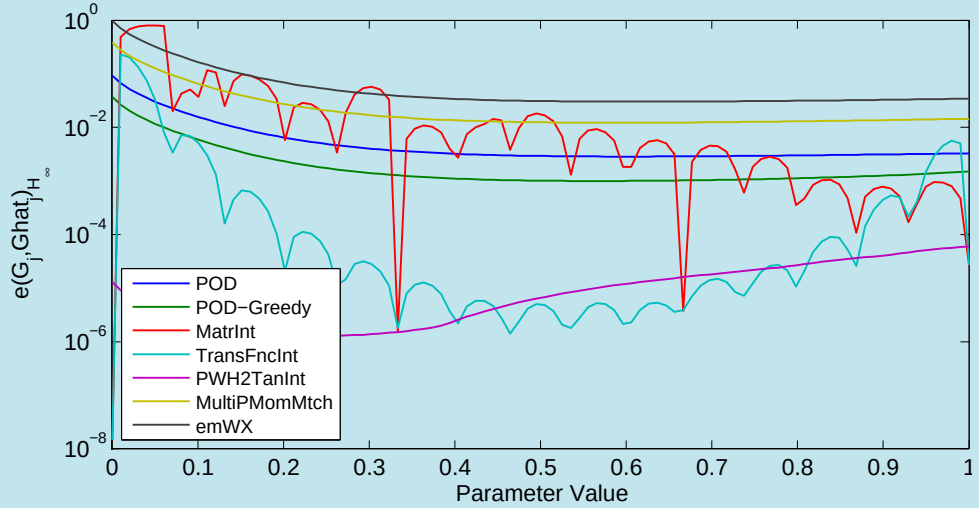
$\mathcal{H}_\infty$ error



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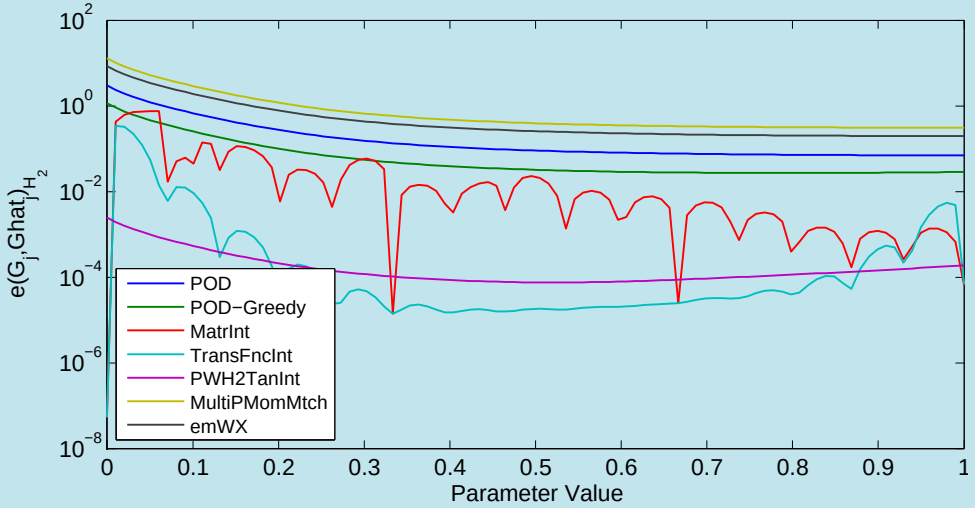
$\mathcal{H}_2$ error



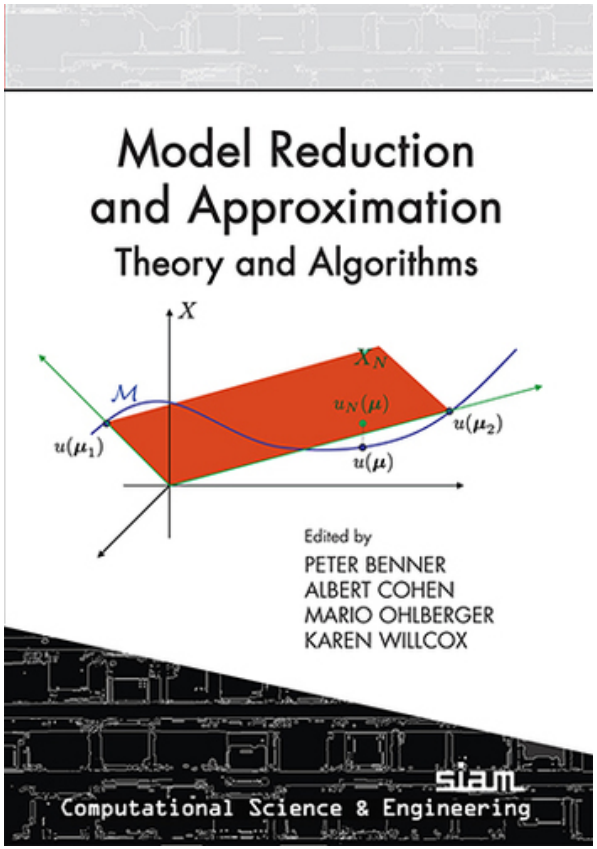
Source: MOR Wiki: http://morwiki.mpi-magdeburg.mpg.de/morwiki/index.php/Anemometer

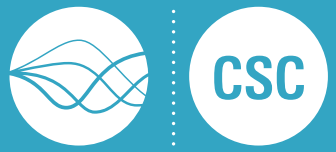
Commercial I

For more details of this comparisons, and other tests, see

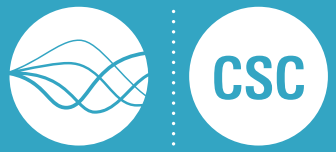
U. Baur, P. Benner, B. Haasdonk, C. Himpe, I. Maier, and M. Ohlberger.
Comparison of Methods for Parametric Model Order Reduction of Unsteady Problems.
In P. Benner, A. Cohen, M. Ohlberger, and K. Willcox (eds.), *Model Reduction and Approximation: Theory and Algorithms.*
SIAM, Philadelphia, PA, 2017.

Chapter 9 in



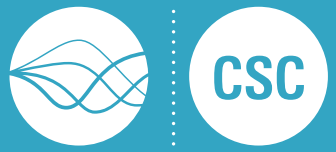


Outline



Conclusions and Outlook

- We have reviewed some of the most popular PMOR methods developed in the last decade, in particular those based on rational interpolation.



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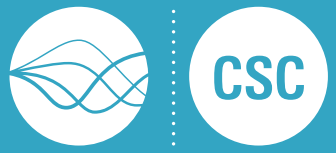

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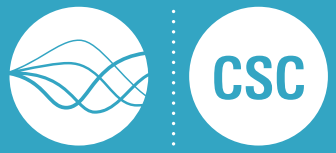
- Employing ideas from reduced-basis method, i.e., greedy sampling, multi-moment matching based PMOR methods can be improved.
- Some ideas from $\mathcal{H}_2$-optimal MOR for non-parametric systems can be extended, but full necessary optimality conditions are still missing.



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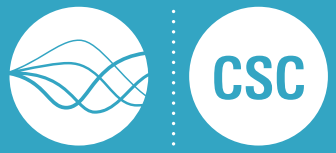
- Some ideas from $\mathcal{H}_2$-optimal MOR for non-parametric systems can be extended, but full necessary optimality conditions are still missing.
- Several extensions to nonlinear systems, but just starting.



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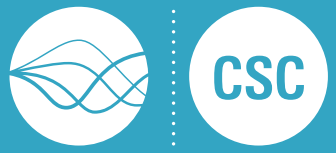

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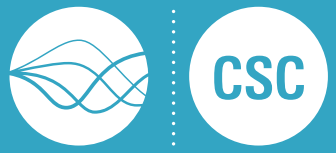
- Most of the methods can be used to significantly accelerate UQ by Monte Carlo or Stochastic Collocation methods!



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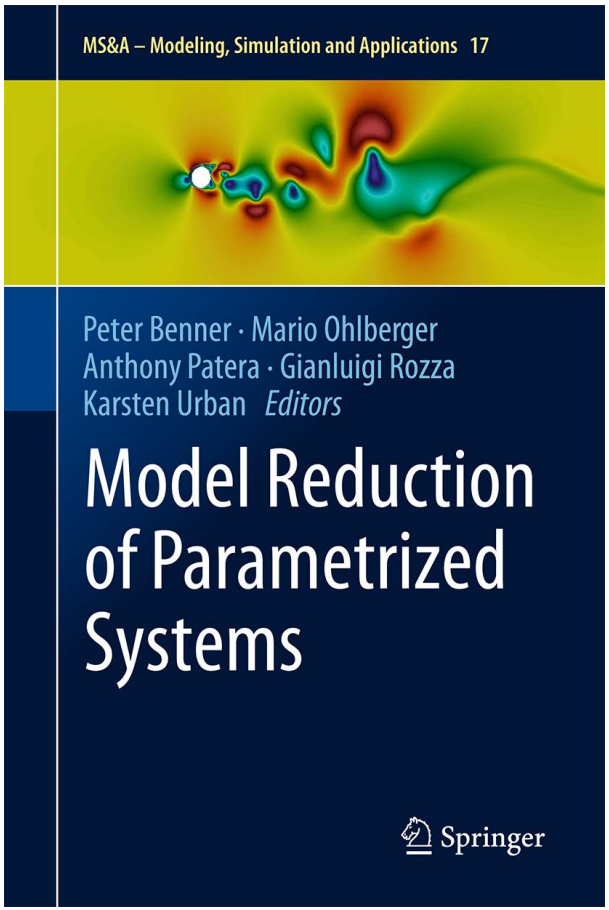

MoRePaS IV — Nantes, April 10–13, 2018.

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