Summer School on Numerical Linear Algebra for Dynamical and High-Dimensional Problems Trogir, October 10–15, 2011

Model Reduction for Linear Dynamical Systems

Peter Benner

Max Planck Institute for Dynamics of Complex Technical Systems Computational Methods in Systems and Control Theory Magdeburg, Germany

http://www.mpi-magdeburg.mpg.de/mpcsc/benner/talks/lecture-MOR.pdf

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- Introduction
- Application Areas
- Motivation
- Model Reduction for Dynamical Systems
- Qualitative and Quantitative Study of the Approximation Error
- Model Reduction by Projection
 - Projection Basics
 - Modal Truncation

Balanced Truncation

- The basic method
- ADI Methods for Lyapunov Equations
- Factored Galerkin-ADI Iteration
- Balancing-Related Model Reduction

Interpolatory Model Reduction

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- Short Introduction
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Microthruster



Final Remarks

Introduction Model Reduction — Abstract Definition

Problem

Given a physical problem with dynamics described by the states $x \in \mathbb{R}^n$, where n is the dimension of the state space.

Because of redundancies, complexity, etc., we want to describe the dynamics of the system using a reduced number of states.

This is the task of model reduction (also: dimension reduction, order reduction).

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- input = output of plant,
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 $\begin{array}{l} \mbox{Modern (LQG-}/\mathcal{H}_{2^{-}}/\mathcal{H}_{\infty}\text{-}) \mbox{ control} \\ \mbox{design: } N \geq n. \end{array}$



Practical controllers require small N ($N \sim 10$, say) due to

- real-time constraints,
- increasing fragility for larger N.

 \implies reduce order of plant (*n*) and/or controller (*N*).



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Application Areas Micro Electronics/Circuit Simulation

- Progressive miniaturization: Moore's Law states that the number of on-chip transistors doubles each 12 (now: 18) months.
- Verification of VLSI/ULSI chip design requires high number of simulations for different input signals.
- Increase in packing density requires modeling of interconncet to ensure that thermic/electro-magnetic effects do not disturb signal transmission.
- Linear systems in micro electronics occur through modified nodal analysis (MNA) for RLC networks, e.g., when
 - decoupling large linear subcircuits,
 - modeling transmission lines (interconnect, powergrid), parasitic effects,
 - modeling pin packages in VLSI chips,
 - modeling circuit elements described by Maxwell's equation using partial element equivalent circuits (PEEC).

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- Resolving complex 3D geometries \Rightarrow millions of degrees of freedom.
- Analysis of elastic deformations requires many simulation runs for varying external forces.

Standard MOR techniques in structural mechanics: modal truncation, combined with Guyan reduction (static condensation) \rightsquigarrow Craig-Bampton method.

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Motivation: Image Compression by Truncated SVD

- A digital image with $n_x \times n_y$ pixels can be represented as matrix $X \in \mathbb{R}^{n_x \times n_y}$, where x_{ij} contains color information of pixel (i, j).
- Memory: $4 \cdot n_x \cdot n_y$ bytes.

Theorem: (Schmidt-Mirsky/Eckart-Young)

Best rank-*r* approximation to $X \in \mathbb{R}^{n_x \times n_y}$ w.r.t. spectral norm:

$$\widehat{X} = \sum_{j=1}^{r} \sigma_j u_j v_j^{T},$$

where $X = U\Sigma V^{T}$ is the singular value decomposition (SVD) of X. The approximation error is $||X - \hat{X}||_2 = \sigma_{r+1}$.

Idea for dimension reduction

Instead of X save $u_1, \ldots, u_r, \sigma_1 v_1, \ldots, \sigma_r v_r$. \rightsquigarrow memory = $4r \times (n_x + n_y)$ bytes.

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Example: Image Compression by Truncated SVD



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• rank r = 50, ≈ 104 kb



Example: Image Compression by Truncated SVD



• rank r = 50, ≈ 104 kb



• rank r = 20, ≈ 42 kb

Rank-20 approximation



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Dimension Reduction via SVD

Example: Gatlinburg

Organizing committee Gatlinburg/Householder Meeting 1964: James H. Wilkinson, Wallace Givens, George Forsythe, Alston Householder, Peter Henrici, Fritz L. Bauer.





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Background: Singular Value Decay

Image data compression via SVD works, if the singular values decay (exponentially).



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Model Reduction for Dynamical Systems

Dynamical Systems

$$\Sigma : \begin{cases} \dot{x}(t) = f(t, x(t), u(t)), & x(t_0) = x_0, \\ y(t) = g(t, x(t), u(t)) \end{cases}$$

with

• states
$$x(t) \in \mathbb{R}^n$$

• inputs
$$u(t) \in \mathbb{R}^m$$

• outputs
$$y(t) \in \mathbb{R}^p$$





Goal:

 $||y - \hat{y}|| < \text{tolerance} \cdot ||u||$ for all admissible input signals



U.

<u>u</u> <u>y</u>



Goal: $\|y - \hat{y}\| < \text{tolerance} \cdot \|u\|$ for all admissible input signals.



Goal:

 $||y - \hat{y}|| < \text{tolerance} \cdot ||u||$ for all admissible input signals. Secondary goal: reconstruct approximation of x from \hat{x} .

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ż	=	f(t, x, u)	=	Ax + Bu,	$A \in \mathbb{R}^{n \times n}$,
y	=	g(t, x, u)	=	Cx + Du,	$C \in \mathbb{R}^{p \times n}$,

State-Space Description for I/O-Relation

 $\mathsf{Variation}\text{-}\mathsf{of}\text{-}\mathsf{constants}\Longrightarrow$

$$\mathcal{S}: u \mapsto y, \quad y(t) = \int_{-\infty}^t C e^{\mathcal{A}(t- au)} \mathcal{B}u(au) \, d au \quad ext{for all } t \in \mathbb{R}$$

 $B \in \mathbb{R}^{n \times m}, \\ D \in \mathbb{R}^{p \times m}.$

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- $S : U \to Y$ is a linear operator between (function) spaces.
- Recall: $A \in \mathbb{R}^{n \times m}$ is a linear operator, $A : \mathbb{R}^m \to \mathbb{R}^n!$
- Basic Idea: use SVD approximation as for matrix A!
- Problem: in general, S does not have a discrete SVD and can therefore not be approximated as in the matrix case!

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Alternative to State-Space Operator: Hankel operator

Instead of

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Linear Systems in Frequency Domain

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Assumptions: $t_0 = 0$, $x_0 = x(0) = 0$.

Laplace Transform / Frequency Domain

Application of Laplace transform

$$\mathcal{L}: x(t) \mapsto x(s) = \int_0^\infty e^{-st} x(t) dt \quad (\Rightarrow \dot{x}(t) \mapsto sx(s))$$

with $s \in \mathbb{C}$ leads to linear system of equations:

$$sx(s) = Ax(s) + Bu(s), \quad y(s) = Cx(s) + Du(s).$$

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Linear Systems in Frequency Domain

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Laplace Transform / Frequency Domain

$$sx(s) = Ax(s) + Bu(s), \quad y(s) = Cx(s) + Du(s)$$

yields I/O-relation in frequency domain:

$$y(s) = \left(\underbrace{C(sI_n - A)^{-1}B + D}_{=:G(s)}\right)u(s) = G(s)u(s).$$

 $G \text{ is the transfer function of } \Sigma, \ G: \mathcal{L}_2^m \to \mathcal{L}_2^p \quad (\mathcal{L}_2:=\mathcal{L}(L_2(-\infty,\infty))).$

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Model Reduction as Approximation Problem

Approximation Problem

Approximate the dynamical system

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by reduced-order system

$$\begin{aligned} \dot{\hat{x}} &= \hat{A}\hat{x} + \hat{B}u, \qquad \hat{A} \in \mathbb{R}^{r \times r}, \quad \hat{B} \in \mathbb{R}^{r \times m}, \\ \dot{\hat{y}} &= \hat{C}\hat{x} + \hat{D}u, \qquad \hat{C} \in \mathbb{R}^{p \times r}, \quad \hat{D} \in \mathbb{R}^{p \times m}. \end{aligned}$$

of order $r \ll n$, such that

$$\|y - \hat{y}\| = \|Gu - \hat{G}u\| \le \|G - \hat{G}\|\|u\| \le ext{tolerance} \cdot \|u\|.$$

 \implies Approximation problem: $\min_{\text{order}(\hat{G}) \leq r} \|G - \hat{G}\|.$

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Model Reduction as Approximation Problem

Approximation Problem

Approximate the dynamical system

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$$\begin{aligned} \dot{\hat{x}} &= \hat{A}\hat{x} + \hat{B}u, \qquad \hat{A} \in \mathbb{R}^{r \times r}, \quad \hat{B} \in \mathbb{R}^{r \times m}, \\ \dot{\hat{y}} &= \hat{C}\hat{x} + \hat{D}u, \qquad \hat{C} \in \mathbb{R}^{p \times r}, \quad \hat{D} \in \mathbb{R}^{p \times m}. \end{aligned}$$

of order $r \ll n$, such that

$$\|y - \hat{y}\| = \|\mathsf{G}u - \hat{\mathsf{G}}u\| \le \|\mathsf{G} - \hat{\mathsf{G}}\|\|u\| \le \mathsf{tolerance} \cdot \|u\|.$$

 \implies Approximation problem: $\min_{\text{order}(\hat{G}) \leq r} \|G - \hat{G}\|$.

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Consider transfer function

$$G(s) = C \left(sI - A \right)^{-1} B + D$$

and input functions $u \in \mathcal{L}_2^m \cong \mathcal{L}_2^m(-\infty,\infty)$, with the 2-norm

$$\|u\|_2^2 := \frac{1}{2\pi} \int_{-\infty}^{\infty} u^*(j\omega) u(j\omega) \, d\omega.$$

Assume A is (asympotically) stable: $\Lambda(A) \subset \mathbb{C}^- := \{z \in \mathbb{C} : re(z) < 0\}.$

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$$\int_{-\infty}^{\infty} y^*(j\omega) y(j\omega) \, d\omega \quad = \quad \int_{-\infty}^{\infty} u^*(j\omega) G^*(j\omega) G(j\omega) u(j\omega) \, d\omega$$

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$$= \int_{-\infty}^{\infty} \|G(j\omega) u(j\omega)\|^2 \, d\omega \le \int_{-\infty}^{\infty} M^2 \|u(j\omega)\|^2 \, d\omega$$

(Here:, ||. || denotes the Euclidian vector or spectral matrix norm.)

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$$\begin{split} \int_{-\infty}^{\infty} y^*(j\omega) y(j\omega) \, d\omega &= \int_{-\infty}^{\infty} u^*(j\omega) G^*(j\omega) G(j\omega) u(j\omega) \, d\omega \\ &= \int_{-\infty}^{\infty} \|G(j\omega) u(j\omega)\|^2 \, d\omega \le \int_{-\infty}^{\infty} M^2 \|u(j\omega)\|^2 \, d\omega \\ &= M^2 \int_{-\infty}^{\infty} u(j\omega)^* u(j\omega) \, d\omega < \infty. \end{split}$$

(Here:, $\|.\|$ denotes the Euclidian vector or spectral matrix norm.)

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$$= M^2 \int_{-\infty}^{\infty} u(j\omega)^* u(j\omega) \, d\omega < \infty.$$

 $\implies y \in L^p_2(-\infty,\infty) \cong \mathcal{L}^p_2.$

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$$\|G\|_{\infty} := \sup_{\|u\|_{2} \neq 0} \frac{\|Gu\|_{2}}{\|u\|_{2}}$$

is well defined. It can be shown that

$$\|G\|_{\infty} := \sup_{\omega \in \mathbb{R}} \|G(\jmath \omega)\| = \sup_{\omega \in \mathbb{R}} \sigma_{max} \left(G(\jmath \omega)\right).$$

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Sketch of proof: $\|G(\jmath\omega)u(\jmath\omega)\| \le \|G(\jmath\omega)\| \|u(\jmath\omega)\| \Rightarrow "\le ".$ Construct u with $\|Gu\|_2 = \sup_{\omega \in \mathbb{R}} \|G(\jmath\omega)\| \|u\|_2.$

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Consider transfer function

$$G(s) = C \left(sI - A \right)^{-1} B + D.$$

Hardy space \mathcal{H}_{∞}

Function space of matrix-/scalar-valued functions that are analytic and bounded in $\mathbb{C}^+.$

The \mathcal{H}_{∞} -norm is

$$\|F\|_{\infty} := \sup_{\mathsf{re}\,s>0} \sigma_{\mathsf{max}}\left(F(s)\right) = \sup_{\omega\in\mathbb{R}} \sigma_{\mathit{max}}\left(F(\jmath\omega)\right).$$

Stable transfer functions are in the Hardy spaces

- \mathcal{H}_{∞} in the SISO case (single-input, single-output, m = p = 1);
- $\mathcal{H}^{p imes m}_{\infty}$ in the MIMO case (multi-input, multi-output, m > 1, p > 1).

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Consider transfer function

$$G(s) = C \left(sI - A \right)^{-1} B + D.$$

Paley-Wiener Theorem (Parseval's equation/Plancherel Theorem)

$$L_2(-\infty,\infty)\cong \mathcal{L}_2, \quad L_2(0,\infty)\cong \mathcal{H}_2$$

Consequently, 2-norms in time and frequency domains coincide!

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\mathcal{H}_{∞} approximation error

Reduced-order model \Rightarrow transfer function $\hat{G}(s) = \hat{C}(sI_r - \hat{A})^{-1}\hat{B} + \hat{D}$. $\|y - \hat{y}\|_2 = \|Gu - \hat{G}\hat{u}\|_2 \le \|G - \hat{G}\|_{\infty} \|u\|_2$.

 \implies compute reduced-order model such that $\|G - \hat{G}\|_{\infty} < to!$ Note: error bound holds in time- and frequency domain due to Paley-Wiener!

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Consider transfer function

$$G(s) = C (sI - A)^{-1} B$$
, i.e. $D = 0$.

Hardy space \mathcal{H}_2

Function space of matrix-/scalar-valued functions that are analytic \mathbb{C}^+ and bounded w.r.t. the $\mathcal{H}_2\text{-norm}$

$$\|F\|_{2} := \left(\sup_{\operatorname{re}\sigma>0}\int_{-\infty}^{\infty}\|F(\sigma+j\omega)\|_{F} d\omega\right)^{\frac{1}{2}}$$
$$= \left(\int_{-\infty}^{\infty}\|F(j\omega)\|_{F} d\omega\right)^{\frac{1}{2}}.$$

Stable transfer functions are in the Hardy spaces

- \mathcal{H}_2 in the SISO case (single-input, single-output, m = p = 1);
- $\mathcal{H}_2^{p \times m}$ in the MIMO case (multi-input, multi-output, m > 1, p > 1).

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$$\|F\|_2 = \left(\int_{-\infty}^{\infty} \|F(j\omega)\|_F \, d\omega\right)^{\frac{1}{2}}$$

 \mathcal{H}_2 approximation error for impulse response $(u(t) = u_0 \delta(t))$

Reduced-order model \Rightarrow transfer function $\hat{G}(s) = \hat{C}(sI_r - \hat{A})^{-1}\hat{B}$.

$$\|y - \hat{y}\|_2 = \|Gu_0\delta - \hat{G}u_0\delta\|_2 \le \|G - \hat{G}\|_2\|u_0\|.$$

 \implies compute reduced-order model such that $\|G - \hat{G}\|_2 < to!!$

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Qualitative and Quantitative Study of the Approximation Error Approximation Problems

\mathcal{H}_{∞} -norm	best approximation problem for given reduced order r in
	general open; balanced truncation yields suboptimal solu-
	tion with computable \mathcal{H}_∞ -norm bound.
\mathcal{H}_2 -norm	necessary conditions for best approximation known; (local)
	optimizer computable with iterative rational Krylov algo-
	rithm (IRKA)
Hankel-norm	optimal Hankel norm approximation (AAK theory).
$\ G\ _H := \sigma_{\max}$	

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Qualitative and Quantitative Study of the Approximation Error Computable error measures

Evaluating system norms is computationally very (sometimes too) expensive.

Other measures

- absolute errors $\|G(\jmath\omega_j) \hat{G}(\jmath\omega_j)\|_2$, $\|G(\jmath\omega_j) \hat{G}(\jmath\omega_j)\|_\infty$ $(j = 1, ..., N_\omega)$; • relative errors $\frac{\|G(\jmath\omega_j) - \hat{G}(\jmath\omega_j)\|_2}{\|G(\jmath\omega_j)\|_2}$, $\frac{\|G(\jmath\omega_j) - \hat{G}(\jmath\omega_j)\|_\infty}{\|G(\jmath\omega_j)\|_\infty}$;
- "eyeball norm", i.e. look at frequency response/Bode (magnitude) plot: for SISO system, log-log plot frequency vs. $|G(j\omega)|$ (or $|G(j\omega) - \hat{G}(j\omega)|$) in decibels, 1 dB $\simeq 20 \log_{10}(\text{value})$.

For MIMO systems, $p \times m$ array of of plots G_{ij} .



	MOR by Projection			
Model _{Goals}	Reduction b	y Projection		

- Automatic generation of compact models.
- Satisfy desired error tolerance for all admissible input signals, i.e., want

 \implies Need computable error bound/estimate!

- Preserve physical properties:
 - stability (poles of G in \mathbb{C}^-),
 - minimum phase (zeroes of G in \mathbb{C}^-),
 - passivity

 $\int_{-\infty}^{t} u(\tau)^{\mathsf{T}} y(\tau) \, d\tau \ge 0 \quad \forall t \in \mathbb{R}, \quad \forall u \in L_2(\mathbb{R}, \mathbb{R}^m).$

	MOR by Projection				
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 - passivity

 $\int_{-\infty}^t u(\tau)^T y(\tau) \, d\tau \ge 0 \quad \forall t \in \mathbb{R}, \quad \forall u \in L_2(\mathbb{R}, \mathbb{R}^m).$

Linear Algebra Basics

Projector

A projector is a matrix $P \in \mathbb{R}^{n \times n}$ with $P^2 = P$. Let $\mathcal{V} = \operatorname{range}(P)$, then P is projector onto \mathcal{V} . On the other hand, if $\{v_1, \ldots, v_r\}$ is a basis of \mathcal{V} and $V = [v_1, \ldots, v_r]$, then $P = V(V^T V)^{-1}V^T$ is a projector onto \mathcal{V} .

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Nodel Reduction by Projection Linear Algebra Basics

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- If $P = P^T$, then P is an orthogonal projector (aka: Galerkin projection), otherwise an oblique projector. (aka: Petrov-Galerkin projection.)
- *P* is the identity operator on \mathcal{V} , i.e., $Pv = v \ \forall v \in \mathcal{V}$.
- I P is the complementary projector onto ker P.
- If V is an A-invariant subspace corresponding to a subset of A's spectrum, then we call P a spectral projector.
- Let $\mathcal{W} \subset \mathbb{R}^n$ be another *r*-dimensional subspace and $W = [w_1, \ldots, w_r]$ be a basis matrix for \mathcal{W} , then $P = V(W^T V)^{-1}W^T$ is an oblique projector onto \mathcal{V} along \mathcal{W} .

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Model F	Reduction b	v Projection		

MOR Methods Based on Projection

Methods:

- Modal Truncation
- Rational Interpolation (Padé-Approximation and (rational) Krylov Subspace Methods)
- Balanced Truncation
- many more...

Joint feature of these methods:

computation of reduced-order model (ROM) by projection!

Model Reduction by Projection

Joint feature of these methods: computation of reduced-order model (ROM) by projection!

Assume trajectory x(t; u) is contained in low-dimensional subspace \mathcal{V} . Thus, use Galerkin or Petrov-Galerkin-type projection of state-space onto \mathcal{V} along complementary subspace \mathcal{W} : $x \approx V W^T x =: \tilde{x}$, where

range
$$(V) = \mathcal{V}$$
, range $(W) = \mathcal{W}$, $W^T V = I_r$.

Then, with $\hat{x} = W^T x$, we obtain $x \approx V \hat{x}$ so that

$$\|x-\tilde{x}\|=\|x-V\hat{x}\|,$$

and the reduced-order model is

 $\hat{A} := W^T A V, \quad \hat{B} := W^T B, \quad \hat{C} := C V, \quad (\hat{D} := D).$
Introduction MOR by Projection Balanced Truncation RatInt Examples Fin.

Model Reduction by Projection

Joint feature of these methods: computation of reduced-order model (ROM) by projection!

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 $\hat{A} := W^T A V, \quad \hat{B} := W^T B, \quad \hat{C} := C V, \quad (\hat{D} := D).$

Important observations:

• The state equation residual satisfies $\dot{\tilde{x}} - A\tilde{x} - Bu \perp \mathcal{W}$, since

$$W^{T}\left(\dot{\tilde{x}} - A\tilde{x} - Bu\right) = W^{T}\left(VW^{T}\dot{x} - AVW^{T}x - Bu\right)$$

Model Reduction by Projection

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$$= \underbrace{W^{T}\dot{x}}_{\dot{\tilde{x}}} - \underbrace{W^{T}AV}_{=\hat{A}}\underbrace{W^{T}x}_{=\hat{x}} - \underbrace{W^{T}B}_{=\hat{B}}u$$

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$$= \dot{\hat{x}} - \hat{A}\hat{x} - \hat{B}u = 0.$$

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IVIOAEI REDUCTION DY Projection MOR Methods Based on Projection

Projection ~ Rational Interpolation

Given the ROM

$$\hat{A} = W^T A V, \quad \hat{B} = W^T B, \quad \hat{C} = C V, \quad (\hat{D} = D),$$

the error transfer function can be written as

$$G(s) - \hat{G}(s) = (C(sI_n - A)^{-1}B + D) - (\hat{C}(sI_n - \hat{A})^{-1}\hat{B} + \hat{D})$$

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= $C((sI_n - A)^{-1} - V(sI_r - \hat{A})^{-1}W^T)B$

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= $C\left((sI_n - A)^{-1} - V(sI_r - \hat{A})^{-1}W^T\right)B$
= $C\left(I_n - \underbrace{V(sI_r - \hat{A})^{-1}W^T(sI_n - A)}_{=:P(s)}\right)(sI_n - A)^{-1}B.$

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IVIOAEI REDUCTION DY Projection MOR Methods Based on Projection

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= $C\left(l_n - \underbrace{V(sl_r - \hat{A})^{-1}W^T(sl_n - A)}_{=:P(s)}\right)(sl_n - A)^{-1}B$

P(s) is a projector onto \mathcal{V} :

range $(P(s)) \subset$ range (V), all matrices have full rank \Rightarrow "=", and $P(s)^2 = V(sI_r - \hat{A})^{-1}W^T(sI_n - A)V(sI_r - \hat{A})^{-1}W^T(sI_n - A)$

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Model Reduction by Projection

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the error transfer function can be written as

$$G(s) - \hat{G}(s) = \left(C(sI_n - A)^{-1}B + D\right) - \left(\hat{C}(sI_n - \hat{A})^{-1}\hat{B} + \hat{D}\right)$$

= $C\left(I_n - \underbrace{V(sI_r - \hat{A})^{-1}W^{T}(sI_n - A)}_{=:P(s)}\right)(sI_n - A)^{-1}B$

P(s) is a projector onto \mathcal{V} :

range $(P(s)) \subset$ range (V), all matrices have full rank \Rightarrow "=", and $P(s)^{2} = V(sl_{r} - \hat{A})^{-1}W^{T}(sl_{n} - A)V(sl_{r} - \hat{A})^{-1}W^{T}(sl_{n} - A)$ $= V(sl_{r} - \hat{A})^{-1}\underbrace{(sl_{r} - \hat{A})(sl_{r} - \hat{A})^{-1}}_{=l_{r}}W^{T}(sl_{n} - A) = P(s).$

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Model Reduction by Projection

Projection ~> Rational Interpolation

Given the ROM

$$\hat{A} = W^T A V, \quad \hat{B} = W^T B, \quad \hat{C} = C V, \quad (\hat{D} = D),$$

the error transfer function can be written as

$$G(s) - \hat{G}(s) = \left(C(sl_n - A)^{-1}B + D\right) - \left(\hat{C}(sl_n - \hat{A})^{-1}\hat{B} + \hat{D}\right)$$

= $C\left(l_n - \underbrace{V(sl_r - \hat{A})^{-1}W^T(sl_n - A)}_{=:P(s)}\right)(sl_n - A)^{-1}B$

 $P(s) \text{ is a projector onto } \mathcal{V} \Longrightarrow$ Given $s_* \in \mathbb{C} \setminus (\Lambda(A) \cup \Lambda(\hat{A})),$ if $(s_*I_n - A)^{-1}B \in \mathcal{V}$, then $(I_n - P(s_*))(s_*I_n - A)^{-1}B = 0,$

hence

$$G(s_*) - \hat{G}(s_*) = 0 \implies G(s_*) = \hat{G}(s_*), \text{ i.e., } \hat{G} \text{ interpolates } G \text{ in } s_*!$$

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Model	Reduction b	y Projection		

MOR Methods Based on Projection

Projection ~> Rational Interpolation

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$$G(s) - \hat{G}(s) = \left(C(sI_n - A)^{-1}B + D\right) - \left(\hat{C}(sI_n - \hat{A})^{-1}\hat{B} + \hat{D}\right) \\ = C\left(I_n - \underbrace{V(sI_r - \hat{A})^{-1}W^{T}(sI_n - A)}_{=:P(s)}\right)(sI_n - A)^{-1}B$$

Analogously, = $C(sI_n - A)^{-1} (I_n - \underbrace{(sI_n - A)V(sI_r - \hat{A})^{-1}W^{\mathsf{T}}}_{=:Q(s)})B.$

 $Q(s)^*$ is a projector onto $\mathcal{W} \Longrightarrow$ Given $s_* \in \mathbb{C} \setminus (\Lambda(A) \cup \Lambda(\hat{A}))$,

if
$$(s_*I_n - A)^{-*}C^T \in W$$
, then $C(s_*I_n - A)^{-1}(I_n - Q(s_*)) = 0$,

hence

$$G(s_*) - \hat{G}(s_*) = 0 \Rightarrow G(s_*) = \hat{G}(s_*), \text{ i.e., } \hat{G} \text{ interpolates } G \text{ in } s_*!$$

Max Planck Institute Magdeburg

Model Reduction by Projection MOR Methods Based on Projection

Theorem

[Grimme '97, Villemagne/Skelton '87]

Given the ROM

$$\hat{A} = W^T A V, \quad \hat{B} = W^T B, \quad \hat{C} = C V, \quad (\hat{D} = D),$$

and $s_* \in \mathbb{C} \setminus (\Lambda(A) \cup \Lambda(\hat{A}))$, if either

•
$$(s_*I_n - A)^{-1}B \in \operatorname{range}(V)$$
, or

•
$$(s_*I_n - A)^{-*}C^T \in \operatorname{range}(W)$$

then the interpolation condition

$$G(s_*)=\hat{G}(s_*).$$

in s* holds.

Note: extension to Hermite interpolation conditions later!

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Basic method:

Assume A is diagonalizable, $T^{-1}AT = D_A$, project state-space onto A-invariant subspace $\mathcal{V} = \operatorname{span}(t_1, \ldots, t_r)$, $v_k = \operatorname{eigenvectors}$ corresp. to "dominant" modes / eigenvalues of A. Then with

 $V = T(:, 1:r) = [t_1, ..., t_r], \quad \tilde{W} = T^{-1}(:, 1:r), \quad W = \tilde{W}(V^*\tilde{W})^{-1},$

reduced-order model is

 $\hat{A} := W^* A V = \operatorname{diag} \{\lambda_1, \dots, \lambda_r\}, \quad \hat{B} := W^* B, \quad \hat{C} = C V$

Also computable by truncation:

$$T^{-1}AT = \begin{bmatrix} \hat{A} \\ A_2 \end{bmatrix}, \quad T^{-1}B = \begin{bmatrix} \hat{B} \\ B_2 \end{bmatrix}, \quad CT = [\hat{C}, C_2], \quad \hat{D} = D.$$

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Properties:

Simple computation for large-scale systems, using, e.g., Krylov subspace methods (Lanczos, Arnoldi), Jacobi-Davidson method.

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Properties:

Error bound:

$$\|G - \hat{G}\|_{\infty} \leq \|C_2\| \|B_2\| \frac{1}{\min_{\lambda \in \Lambda(A_2)} |\operatorname{Re}(\lambda)|}$$

Proof:

$$G(s) = C(sI - A)^{-1}B + D = CTT^{-1}(sI - A)^{-1}TT^{-1}B + D$$

= $CT(sI - T^{-1}AT)^{-1}T^{-1}B + D$
= $[\hat{C}, C_2] \begin{bmatrix} (sI_r - \hat{A})^{-1} \\ (sI_{n-r} - A_2)^{-1} \end{bmatrix} \begin{bmatrix} \hat{B} \\ B_2 \end{bmatrix} + D$
= $\hat{G}(s) + C_2(sI_{n-r} - A_2)^{-1}B_2,$

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Proof:

$$G(s) = \hat{G}(s) + C_2(sI_{n-r} - A_2)^{-1}B_2,$$

observing that $\|G - \hat{G}\|_{\infty} = \sup_{\omega \in \mathbb{R}} \sigma_{\max}(C_2(\jmath \omega I_{n-r} - A_2)^{-1}B_2)$, and

$$C_2(\jmath\omega I_{n-r}-A_2)^{-1}B_2=C_2 {
m diag}\left(rac{1}{\jmath\omega-\lambda_{r+1}},\ldots,rac{1}{\jmath\omega-\lambda_n}
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Difficulties:

- Eigenvalues contain only limited system information.
- Dominance measures are difficult to compute. ([LITZ '79] use Jordan canoncial form; otherwise merely heuristic criteria, e.g., [VARGA '95]. Recent improvement: dominant pole algorithm.)
- Error bound not computable for really large-scale problems.

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Modal Truncation Example

BEAM, SISO system from SLICOT Benchmark Collection for Model Reduction, n = 348, m = p = 1, reduced using 13 dominant complex conjugate eigenpairs, error bound yields $\|G - \hat{G}\|_{\infty} \le 1.21 \cdot 10^3$



MATLAB[®] demo.

Coffee break!

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Basic principle:

 A system Σ, realized by (A, B, C, D), is called balanced, if the Gramians, i.e., solutions P, Q of the Lyapunov equations

$$AP + PA^{\mathsf{T}} + BB^{\mathsf{T}} = 0, \qquad A^{\mathsf{T}}Q + QA + C^{\mathsf{T}}C = 0,$$

satisfy: $P = Q = \operatorname{diag}(\sigma_1, \dots, \sigma_n)$ with $\sigma_1 \ge \sigma_2 \ge \dots \ge \sigma_n > 0$.

• $\Lambda(PQ)^{\frac{1}{2}} = \{\sigma_1, \dots, \sigma_n\}$ are the Hankel singular values (HSVs) of Σ .

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$$y(t) = \mathcal{H}u(t) = \int_{-\infty}^{0} Ce^{A(t-\tau)} Bu(\tau) d\tau$$

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$$\mathcal{H}^*\mathcal{H}u(t) = B^T e^{-A^T t} Qz \doteq \sigma^2 u(t).$$

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$$\frac{1}{\sigma^2} \int_{-\infty}^0 e^{-A\tau} B \frac{1}{\sigma^2} B^T e^{-A^T \tau} Q z d\tau$$

$$= \frac{1}{\sigma^2} \int_{-\infty}^{\infty} e^{At} BB^T e^{A^T t} dt Qz$$
$$= \frac{1}{\sigma^2} \underbrace{\int_{0}^{\infty} e^{At} BB^T e^{A^T t} dt}_{=0} Qz$$

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z

$$= \int_{-\infty}^{0} e^{-A\tau} B \frac{1}{\sigma^2} B^{T} e^{-A^{T}\tau} Qz d\tau$$
$$= \frac{1}{\sigma^2} \underbrace{\int_{0}^{\infty} e^{At} BB^{T} e^{A^{T}t} dt}_{0} Qz$$

$$\equiv P$$

$$= \frac{1}{\sigma^2} PQz$$

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= $\frac{1}{\sigma^2}PQz$

$$= \int_{-\infty}^{0} e^{-A\tau} B \frac{1}{\sigma^2} B^T e^{-A^T \tau} Qz d\tau$$
$$= \frac{1}{\sigma^2} \int_{-\infty}^{\infty} e^{At} B B^T e^{A^T t} dt Qz$$

$$= \frac{1}{\sigma^2} \underbrace{\int_0^{\sigma^2} e^{i x BB + e^{i x - dt}}}_{\equiv P} Qz$$

$$\iff PQz = \sigma^2 z.$$

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- $\Lambda(PQ)^{\frac{1}{2}} = \{\sigma_1, \ldots, \sigma_n\}$ are the Hankel singular values (HSVs) of Σ .
- Compute balanced realization of the system via state-space transformation

$$\mathcal{T}: (A, B, C, D) \mapsto (TAT^{-1}, TB, CT^{-1}, D)$$

$$= \left(\begin{bmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{bmatrix}, \begin{bmatrix} B_1 \\ B_2 \end{bmatrix}, \begin{bmatrix} C_1 & C_2 \end{bmatrix}, D \right)$$
Fruncation $\rightsquigarrow (\hat{A}, \hat{B}, \hat{C}, \hat{D}) := (A_{11}, B_1, C_1, D).$

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Basic principle:

 A system Σ, realized by (A, B, C, D), is called balanced, if the Gramians, i.e., solutions P, Q of the Lyapunov equations

$$AP + PA^T + BB^T = 0, \qquad A^TQ + QA + C^TC = 0,$$

satisfy: $P = Q = \operatorname{diag}(\sigma_1, \ldots, \sigma_n)$ with $\sigma_1 \ge \sigma_2 \ge \ldots \ge \sigma_n > 0$.

- $\Lambda(PQ)^{\frac{1}{2}} = \{\sigma_1, \ldots, \sigma_n\}$ are the Hankel singular values (HSVs) of Σ .
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Motivation:

HSVs are system invariants: they are preserved under $\mathcal{T} : (A, B, C, D) \mapsto (TAT^{-1}, TB, CT^{-1}, D)$:

in transformed coordinates, the Gramians satisfy

$$(TAT^{-1})(TPT^{T}) + (TPT^{T})(TAT^{-1})^{T} + (TB)(TB)^{T} = 0$$

 $(TAT^{-1})^{T}(T^{-T}QT^{-1}) + (T^{-T}QT^{-1})(TAT^{-1}) + (CT^{-1})^{T}(CT^{-1}) = 0$

 $\Rightarrow (TPT^{T})(T^{-T}QT^{-1}) = TPQT^{-1},$

hence $\Lambda(PQ) = \Lambda((TPT^{T})(T^{-T}QT^{-1})).$

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Motivation:

HSVs are system invariants: they are preserved under $\mathcal{T} : (A, B, C, D) \mapsto (TAT^{-1}, TB, CT^{-1}, D)$

HSVs determine the energy transfer given by the Hankel map

$$\mathcal{H}: L_2(-\infty, 0) \mapsto L_2(0, \infty): u_- \mapsto y_+.$$

In balanced coordinates ... energy transfer from u_{-} to y_{+} :

$$E := \sup_{u \in L_2(-\infty,0] \atop x(0)=x_0} \frac{\int_{0}^{\infty} y(t)^T y(t) dt}{\int_{-\infty}^{0} u(t)^T u(t) dt} = \frac{1}{\|x_0\|_2} \sum_{j=1}^{n} \sigma_j^2 x_{0,j}^2$$

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 \implies Truncate states corresponding to "small" HSVs \implies complete analogy to best approximation via SVD!







Implementation: SR Method

• Compute (Cholesky) factors of the Gramians, $P = S^T S$, $Q = R^T R$.

• Compute SVD $SR^{T} = \begin{bmatrix} U_1, U_2 \end{bmatrix} \begin{bmatrix} \Sigma_1 \\ \Sigma_2 \end{bmatrix} \begin{bmatrix} V_1^{T} \\ V_2^{T} \end{bmatrix}.$

Solution ROM is $(W^T AV, W^T B, CV, D)$, where

$$W = R^T V_1 \Sigma_1^{-\frac{1}{2}}, \qquad V = S^T U_1 \Sigma_1^{-\frac{1}{2}}$$

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$$= \Sigma_{1}^{-\frac{1}{2}}[I_{r}, 0] \begin{bmatrix} \Sigma_{1} \\ \Sigma_{2} \end{bmatrix} \begin{bmatrix} I_{r} \\ 0 \end{bmatrix} \Sigma_{1}^{-\frac{1}{2}}$$

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 $\implies VW^{T}$ is an oblique projector, hence balanced truncation is a Petrov-Galerkin projection method.

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Properties:

- Reduced-order model is stable with HSVs $\sigma_1, \ldots, \sigma_r$.
- Adaptive choice of *r* via computable error bound:

$$||y - \hat{y}||_2 \le \left(2\sum_{k=r+1}^n \sigma_k\right) ||u||_2.$$

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"New" algorithmic ideas from numerical linear algebra:

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Properties:

General misconception: complexity $O(n^3)$ – true for several implementations! (e.g., MATLAB, SLICOT).

- "New" algorithmic ideas from numerical linear algebra:
- Instead of Gramians P, Qcompute $S, R \in \mathbb{R}^{n \times k}$, $k \ll n$, such that

$$P \approx SS^T$$
, $Q \approx RR^T$.

 Compute S, R with problem-specific Lyapunov solvers of "low" complexity directly.



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"New" algorithmic ideas from numerical linear algebra:

Sparse Balanced Truncation:

- − Sparse implementation using sparse Lyapunov solver (→ADI+MUMPS/SuperLU).
- Complexity $\mathcal{O}(n(k^2 + r^2))$.
- Software:
 - + MATLAB toolbox LyaPack (PENZL 1999),
 - + Software library M.E.S.S.^a in C/MATLAB [B./SAAK/KÖHLER].

^aMatrix Equation Sparse Solvers

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ADI Methods for Lyapunov Equations Background

Recall Peaceman Rachford ADI:

Consider Au = s where $A \in \mathbb{R}^{n \times n}$ spd, $s \in \mathbb{R}^n$. ADI Iteration Idea: Decompose A = H + V with $H, V \in \mathbb{R}^{n \times n}$ such that

(H + pI)v = r(V + pI)w = t

can be solved easily/efficiently.

ADI Iteration

If $H, V \text{ spd} \Rightarrow \exists p_k, k = 1, 2, \dots$ such that

$$u_{0} = 0$$

(H + p_{k}l)u_{k-\frac{1}{2}} = (p_{k}l - V)u_{k-1} + s
(V + p_{k}l)u_{k} = (p_{k}l - H)u_{k-\frac{1}{2}} + s

converges to $u \in \mathbb{R}^n$ solving Au = s.

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ADI Methods for Lyapunov Equations

The Lyapunov operator

$$\mathcal{L}: P \mapsto AX + XA^T$$

can be decomposed into the linear operators

 $\mathcal{L}_H: X \mapsto AX, \qquad \mathcal{L}_V: X \mapsto XA^T.$

In analogy to the standard ADI method we find the



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ADI N	Methods for I	_yapunov Equat	ions	

Low-Rank ADI

Consider
$$AX + XA^T = -BB^T$$
 for stable $A, B \in \mathbb{R}^{n \times m}$ with $m \ll n$.

ADI iteration for the Lyapunov equation

[Wachspress '95]

For $k = 1, \ldots, k_{\max}$

$$\begin{array}{rcl} X_0 &= & 0 \\ (A+p_k I) X_{k-\frac{1}{2}} &= & -BB^T - X_{k-1} (A^T - p_k I) \\ (A+p_k I) X_k^T &= & -BB^T - X_{k-\frac{1}{2}}^T (A^T - p_k I) \end{array}$$

Rewrite as one step iteration and factorize $X_k = Z_k Z_k^T$, $k = 0, \ldots, k_{max}$

$$Z_{0}Z_{0}^{T} = 0$$

$$Z_{k}Z_{k}^{T} = -2p_{k}(A + p_{k}I)^{-1}BB^{T}(A + p_{k}I)^{-T} + (A + p_{k}I)^{-1}(A - p_{k}I)Z_{k-1}Z_{k-1}^{T}(A - p_{k}I)^{T}(A + p_{k}I)^{-T}$$

... ~> low-rank Cholesky factor ADI

[PENZL '97/'00, LI/WHITE '99/'02, B./LI/PENZL '99/'08, GUGERCIN/SORENSEN/ANTOULAS '03

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 $\ldots \rightsquigarrow$ low-rank Cholesky factor ADI

[PENZL '97/'00, LI/WHITE '99/'02, B./LI/PENZL '99/'08, GUGERCIN/SORENSEN/ANTOULAS '03

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[PENZL '97/'00, LI/WHITE '99/'02, B./LI/PENZL '99/'08, GUGERCIN/SORENSEN/ANTOULAS '03]

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Balanced Truncation ADI Methods for Lyapunov Equations

$$Z_{k} = [\sqrt{-2p_{k}}(A + p_{k}I)^{-1}B, \ (A + p_{k}I)^{-1}(A - p_{k}I)Z_{k-1}]$$

[PENZL '00]

Observing that $(A - p_i I)$, $(A + p_k I)^{-1}$ commute, we rewrite $Z_{k_{max}}$ as

$$Z_{k_{\max}} = [z_{k_{\max}}, P_{k_{\max}-1}z_{k_{\max}}, P_{k_{\max}-2}(P_{k_{\max}-1}z_{k_{\max}}), \dots, P_1(P_2 \cdots P_{k_{\max}-1}z_{k_{\max}})],$$

where

$$z_{k_{\max}} = \sqrt{-2p_{k_{\max}}}(A + p_{k_{\max}}I)^{-1}B$$

and

$$P_i := \frac{\sqrt{-2p_i}}{\sqrt{-2p_{i+1}}} \left[I - (p_i + p_{i+1})(A + p_i I)^{-1} \right].$$

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Balanced Truncation ADI Methods for Lyapunov Equations

$$Z_{k} = \left[\sqrt{-2p_{k}}(A + p_{k}I)^{-1}B, \ (A + p_{k}I)^{-1}(A - p_{k}I)Z_{k-1}\right]$$

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[LI/WHITE '02

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ADI Methods for Lyapunov Equations Lyapunov equation $0 = AX + XA^T + BB^T$.

Algorithm [Penzl '97/'00, LI/WHITE '99/'02, B. 04, B./LI/PENZL '99/'08]

$$V_{1} \leftarrow \sqrt{-2 \operatorname{re} p_{1}} (A + p_{1}I)^{-1}B, \quad Z_{1} \leftarrow V_{1}$$

FOR $k = 2, 3, ...$
$$V_{k} \leftarrow \sqrt{\frac{\operatorname{re} p_{k}}{\operatorname{re} p_{k-1}}} (V_{k-1} - (p_{k} + \overline{p_{k-1}})(A + p_{k}I)^{-1}V_{k-1})$$
$$Z_{k} \leftarrow [Z_{k-1} \quad V_{k}]$$
$$Z_{k} \leftarrow \operatorname{rrlq}(Z_{k}, \tau) \quad \text{column compression}$$

At convergence, $Z_{k_{\text{max}}} Z_{k_{\text{max}}}^{T} \approx X$, where (without column compression)

$$Z_{k_{\max}} = \begin{bmatrix} V_1 & \dots & V_{k_{\max}} \end{bmatrix}, \quad V_k = \begin{bmatrix} \in \mathbb{C}^{n \times m}. \end{bmatrix}$$

Note: Implementation in real arithmetic possible by combining two steps [B./Li/Penzl '99/'08] or using new idea employing the relation of 2 consecutive complex factors [B./Kürschner/Saak '11].

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Numerical Results for ADI Optimal Cooling of Steel Profiles

 Mathematical model: boundary control for linearized 2D heat equation.

$$c \cdot \rho \frac{\partial}{\partial t} x = \lambda \Delta x, \quad \xi \in \Omega$$

$$\lambda \frac{\partial}{\partial n} x = \kappa (u_k - x), \quad \xi \in \Gamma_k, \ 1 \le k \le 7$$

$$\frac{\partial}{\partial n} x = 0, \qquad \xi \in \Gamma_7.$$

$$\implies m = 7, p = 6.$$

- FEM Discretization, different models for initial mesh (n = 371),
 1, 2, 3, 4 steps of mesh refinement ⇒
 - n = 1357, 5177, 20209, 79841.



Source: Physical model: courtesy of Mannesmann/Demag.

Math. model: TRÖLTZSCH/UNGER 1999/2001, PENZL 1999, SAAK 2003.

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Numerical Results for ADI Optimal Cooling of Steel Profiles

• Solve dual Lyapunov equations needed for balanced truncation, i.e.,

 $APM^{T} + MPA^{T} + BB^{T} = 0, \quad A^{T}QM + M^{T}QA + C^{T}C = 0,$

for 79,841.

- 25 shifts chosen by Penzl heuristic from 50/25 Ritz values of A of largest/smallest magnitude, no column compression performed.
- New version in M.E.S.S. requires no factorization of mass matrix!
- Computations done on Core2Duo at 2.8GHz with 3GB RAM and 32Bit-MATLAB.



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Numerica	al Results fo	or ADI			
Scaling / Mesh	Independence		Computations	by Martin Köh	ler '10

- A ∈ ℝ^{n×n} ≡ FDM matrix for 2D heat equation on [0, 1]² (LYAPACK benchmark demo_l1, m = 1).
- 16 shifts chosen by Penzl heuristic from 50/25 Ritz values of A of largest/smallest magnitude.
- Computations using 2 dual core Intel Xeon 5160 with 16 GB RAM.

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Numerica	I Results fo	r ADI		

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n	M.E.S.S. (C)	LyaPack	M.E.S.S. (MATLAB)		
100	0.023	0.124	0.158		
625	0.042	0.104	0.227		
2,500	0.159	0.702	0.989		
10,000	0.965	6.22	5.644		
40,000	11.09	71.48	34.55		
90,000	34.67	418.5	90.49		
160,000	109.3	out of memory	219.9		
250,000	193.7	out of memory	403.8		
562,500	930.1	out of memory	1216.7		
1,000,000	2220.0	out of memory	2428.6		

CPU Times



- A ∈ ℝ^{n×n} ≡ FDM matrix for 2D heat equation on [0, 1]² (LYAPACK benchmark demo_l1, m = 1).
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Factored Galerkin-ADI Iteration Lyapunov equation $0 = AX + XA^T + BB^T$

Projection-based methods for Lyapunov equations with $A + A^T < 0$:

Ocmpute orthonormal basis range (Z), Z ∈ ℝ^{n×r}, for subspace Z ⊂ ℝⁿ, dim Z = r.

2 Set
$$\hat{A} := Z^T A Z$$
, $\hat{B} := Z^T B$

- **③** Solve small-size Lyapunov equation $\hat{A}\hat{X} + \hat{X}\hat{A}^{T} + \hat{B}\hat{B}^{T} = 0$.
- Use $X \approx Z \hat{X} Z^T$.

Examples:

• Krylov subspace methods, i.e., for m = 1:

$$\mathcal{Z} = \mathcal{K}(A, B, r) = \operatorname{span}\{B, AB, A^2B, \dots, A^{r-1}B\}$$

[SAAD '90, JAIMOUKHA/KASENALLY '94, JBILOU '02-'08].

• K-PIK [Simoncini '07],

$$\mathcal{Z} = \mathcal{K}(A, B, r) \cup \mathcal{K}(A^{-1}, B, r).$$

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$$\hat{A} := Z^T A Z$$
, $\hat{B} := Z^T B$

- **③** Solve small-size Lyapunov equation $\hat{A}\hat{X} + \hat{X}\hat{A}^{T} + \hat{B}\hat{B}^{T} = 0$.
- Use $X \approx Z \hat{X} Z^T$.

Examples:

• Krylov subspace methods, i.e., for m = 1:

$$\mathcal{Z} = \mathcal{K}(A, B, r) = \operatorname{span}\{B, AB, A^2B, \dots, A^{r-1}B\}$$

[SAAD '90, JAIMOUKHA/KASENALLY '94, JBILOU '02-'08].

• K-PIK [Simoncini '07],

$$\mathcal{Z} = \mathcal{K}(A, B, r) \cup \mathcal{K}(A^{-1}, B, r).$$

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Factored Galerkin-ADI Iteration Lyapunov equation $0 = AX + XA^T + BB^T$

Projection-based methods for Lyapunov equations with $A + A^T < 0$:

- Compute orthonormal basis range (Z), Z ∈ ℝ^{n×r}, for subspace Z ⊂ ℝⁿ, dim Z = r.
- **③** Solve small-size Lyapunov equation $\hat{A}\hat{X} + \hat{X}\hat{A}^{T} + \hat{B}\hat{B}^{T} = 0$.
- Use $X \approx Z \hat{X} Z^T$.

Examples:

• ADI subspace [B./R.-C. LI/TRUHAR '08]:

$$\mathcal{Z} = \operatorname{colspan} \left[\begin{array}{cc} V_1, & \dots, & V_r \end{array} \right].$$

Note:

- ADI subspace is rational Krylov subspace [J.-R. LI/WHITE '02].
- Similar approach: ADI-preconditioned global Arnoldi method [JBILOU '08].



Factored Galerkin-ADI Iteration

FEM semi-discretized control problem for parabolic PDE:

- optimal cooling of rail profiles,
- n = 20, 209, m = 7, p = 6.



CPU times: 80s (projection every 5th ADI step) vs. 94s (no projection).



Factored Galerkin-ADI Iteration

FEM semi-discretized control problem for parabolic PDE:

- optimal cooling of rail profiles,
- n = 20, 209, m = 7, p = 6.



CPU times: 368s (projection every 5th ADI step) vs. 1207s (no projection).



Factored Galerkin-ADI Iteration Numerical examples for Galerkin-ADI: optimal cooling of rail profiles, n = 79,841.

MESS w/o Galerkin projection and column compression



MESS with Galerkin projection and column compression



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Balanced Truncation Numerical example for BT: Optimal Cooling of Steel Profiles



- BT model computed with sign function method,
- MT w/o static condensation, same order as BT model.

Balanced Truncation Numerical example for BT: Optimal Cooling of Steel Profiles



- BT model computed with sign function method,
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Balanced Truncation Numerical example for BT: Microgyroscope (Butterfly Gyro)



- By applying AC voltage to electrodes, wings are forced to vibrate in anti-phase in wafer plane.
- Coriolis forces induce motion of wings out of wafer plane yielding sensor data.

- Vibrating micro-mechanical gyroscope for inertial navigation.
- Rotational position sensor.



Source: The Oberwolfach Benchmark Collection http://www.intek.de/simulation/benchmark Courtesy of D. Billger (Imego Institute, Göteborg), Saab Bofors Dynamics AB.

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Balanced Iruncation

Numerical example for BT: Microgyroscope (Butterfly Gyro)

- FEM discretization of structure dynamical model using quadratic tetrahedral elements (ANSYS-SOLID187)
 - → n = 34,722, m = 1, p = 12.
- Reduced model computed using SPARED, r = 30.

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Balanced Truncation

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Basic Principle

Given positive semidefinite matrices $P = S^T S$, $Q = R^T R$, compute balancing state-space transformation so that

$$P = Q = \operatorname{diag}(\sigma_1, \ldots, \sigma_n) = \Sigma, \quad \sigma_1 \ge \ldots \ge \sigma_n \ge 0,$$

and truncate corresponding realization at size r with $\sigma_r > \sigma_{r+1}$.

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Classical Balanced Truncation (BT)

Mullis/Roberts '76, Moore '81

- P =controllability Gramian of system given by (A, B, C, D).
- Q = observability Gramian of system given by (A, B, C, D).
- P, Q solve dual Lyapunov equations

$$AP + PA^{T} + BB^{T} = 0, \qquad A^{T}Q + QA + C^{T}C = 0.$$

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LQG Balanced Truncation (LQGBT) [JONCKHEERE/SILVERMAN '83]

- P/Q = controllability/observability Gramian of closed-loop system based on LQG compensator.
- P, Q solve dual algebraic Riccati equations (AREs)

$$0 = AP + PA^{T} - PC^{T}CP + B^{T}B,$$

$$0 = A^{T}Q + QA - QBB^{T}Q + C^{T}C.$$

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Basic Principle

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$$P = Q = \operatorname{diag}(\sigma_1, \ldots, \sigma_n) = \Sigma, \quad \sigma_1 \ge \ldots \ge \sigma_n \ge 0,$$

and truncate corresponding realization at size r with $\sigma_r > \sigma_{r+1}$.

Balanced Stochastic Truncation (BST) [Desai/Pal '84, Green '88]

- P = controllability Gramian of system given by (A, B, C, D), i.e., solution of Lyapunov equation $AP + PA^T + BB^T = 0$.
- Q = observability Gramian of right spectral factor of power spectrum of system given by (A, B, C, D), i.e., solution of ARE

$$\hat{A}^{\mathsf{T}}Q + Q\hat{A} + QB_{W}(DD^{\mathsf{T}})^{-1}B_{W}^{\mathsf{T}}Q + C^{\mathsf{T}}(DD^{\mathsf{T}})^{-1}C = 0,$$

where $\hat{A} := A - B_W (DD^T)^{-1} C$, $B_W := BD^T + PC^T$.

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Basic Principle

Given positive semidefinite matrices $P = S^T S$, $Q = R^T R$, compute balancing state-space transformation so that

$$P = Q = \operatorname{diag}(\sigma_1, \ldots, \sigma_n) = \Sigma, \quad \sigma_1 \ge \ldots \ge \sigma_n \ge 0,$$

and truncate corresponding realization at size r with $\sigma_r > \sigma_{r+1}$.

Positive-Real Balanced Truncation (PRBT)

- Based on positive-real equations, related to positive real (Kalman-Yakubovich-Popov-Anderson) lemma.
- P, Q solve dual AREs

$$0 = \overline{A}P + P\overline{A}^{T} + PC^{T}\overline{R}^{-1}CP + B\overline{R}^{-1}B^{T},$$

$$0 = \overline{A}^{T}Q + Q\overline{A} + QB\overline{R}^{-1}B^{T}Q + C^{T}\overline{R}^{-1}C,$$

where $\bar{R} = D + D^T$, $\bar{A} = A - B\bar{R}^{-1}C$.

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Basic Principle

Given positive semidefinite matrices $P = S^T S$, $Q = R^T R$, compute balancing state-space transformation so that

$$P = Q = \operatorname{diag}(\sigma_1, \ldots, \sigma_n) = \Sigma, \quad \sigma_1 \ge \ldots \ge \sigma_n \ge 0,$$

and truncate corresponding realization at size r with $\sigma_r > \sigma_{r+1}$.

Other Balancing-Based Methods

- Bounded-real balanced truncation (BRBT) based on bounded real lemma [OPDENACKER/JONCKHEERE '88];
- H_{∞} balanced truncation (HinfBT) closed-loop balancing based on H_{∞} compensator [MUSTAFA/GLOVER '91].

Both approaches require solution of dual AREs.

• Frequency-weighted versions of the above approaches.

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Balancing-Related Model Reduction Properties

- Guaranteed preservation of physical properties like
 - stability (all),
 - passivity (PRBT),
 - minimum phase (BST).
- Computable error bounds, e.g.,

$$\begin{split} \mathsf{BT:} & \|G - G_r\|_{\infty} &\leq 2 \; \sum_{j=r+1}^{n} \sigma_j^{BT}, \\ \mathsf{LQGBT:} & \|G - G_r\|_{\infty} &\leq \; 2 \sum_{j=r+1}^{n} \frac{\sigma_j^{LQG}}{\sqrt{1 + (\sigma_j^{LQG})^2}} \\ \mathsf{BST:} & \|G - G_r\|_{\infty} &\leq \left(\prod_{j=r+1}^{n} \frac{1 + \sigma_j^{BST}}{1 - \sigma_j^{BST}} - 1\right) \|G\|_{\infty}, \end{split}$$

- Can be combined with singular perturbation approximation for steady-state performance.
- Computations can be modularized.

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Idea:

Consider

$$\dot{x} = Ax + Bu, \quad y = Cx$$

with transfer function $G(s) = C(sI_n - A)^{-1}B$.

• For $s_0 \notin \Lambda(A)$:

$$G(s) = C \left(I - (s - s_0)(s_0 I_n - A)^{-1} \right)^{-1} (s_0 I_n - A)^{-1} B$$

= $m_0 + m_1(s - s_0) + m_2(s - s_0)^2 + \dots$

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= $C(I - (s - s_0)(s_0I_n - A)^{-1})^{-1}(s_0I_n - A)^{-1}B$
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- For
$$s_0 = 0$$
: $m_j := C(A^{-1})^j B$ = moments.
- For $s_0 = \infty$: $m_j := CA^{j-1}B$ = Markov parameters.

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= $m_0 + m_1(s - s_0) + m_2(s - s_0)^2 + \dots$

• As reduced-order model use *r*th Padé approximant \hat{G} to *G*:

$$G(s) = \hat{G}(s) + \mathcal{O}((s-s_0)^{2r}),$$

i.e., $m_j = \widehat{m}_j$ for $j = 0, \ldots, 2r - 1$

 \rightsquigarrow moment matching if $s_0 < \infty$,

 \rightsquigarrow partial realization if $s_0 = \infty$.

Padé-via-Lanczos Method (PVL)

• Moments need not be computed explicitly; moment matching is equivalent to projecting state-space onto

 $\mathcal{V} = \operatorname{span}(\tilde{B}, \tilde{A}\tilde{B}, \dots, \tilde{A}^{r-1}\tilde{B}) =: \mathcal{K}(\tilde{A}, \tilde{B}, r)$

(where $\tilde{A} = (s_0 I_n - A)^{-1}$, $\tilde{B} = (s_0 I_n - A)^{-1}B$) along

$$\mathcal{W} = \operatorname{span}(\mathcal{C}^{\mathsf{T}}, \tilde{\mathcal{A}}^* \mathcal{C}^{\mathsf{T}}, \dots, (\tilde{\mathcal{A}}^*)^{r-1} \mathcal{C}^{\mathsf{T}}) =: \mathcal{K}(\tilde{\mathcal{A}}^*, \mathcal{C}^{\mathsf{T}}, r).$$

• Computation via unsymmetric Lanczos method, yields system matrices of reduced-order model as by-product.

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• Computation via unsymmetric Lanczos method, yields system matrices of reduced-order model as by-product.

Remark: Arnoldi (PRIMA) yields only $G(s) = \hat{G}(s) + O((s - s_0)^r)$.

Padé-via-Lanczos Method (PVL)

- Computable error estimates/bounds for $\|y \hat{y}\|_2$ often very pessimistic or expensive to evaluate.
- Mostly heuristic criteria for choice of expansion points. Optimal choice for second-order systems with proportional/Rayleigh damping (BEATTIE/GUGERCIN '05).
- Good approximation quality only locally.
- Preservation of physical properties only in special cases; usually requires post processing which (partially) destroys moment matching properties.

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Internolatory Model		I Reduction		
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Computation of reduced-order model by projection

Given an LTI system $\dot{x} = Ax + Bu$, y = Cx with transfer function $G(s) = C(sI_n - A)^{-1}B$, a reduced-order model is obtained using projection approach with $V, W \in \mathbb{R}^{n \times r}$ and $W^T V = I_r$ by computing

$$\hat{A} = W^T A V, \quad \hat{B} = W^T B, \quad \hat{C} = C V.$$

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

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

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

Rational Interpolation/Moment-Matching

Choose V, W such that

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

and

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

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Theorem (simplified) [GRIMME '97, VILLEMAGNE/SKELTON '87]

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$$\begin{array}{ll} \operatorname{span}\left\{(s_1I_n-A)^{-1}B,\ldots,(s_kI_n-A)^{-1}B\right\} &\subset & \operatorname{Ran}(V), \\ \operatorname{span}\left\{(s_1I_n-A)^{-T}C^T,\ldots,(s_kI_n-A)^{-T}C^T\right\} &\subset & \operatorname{Ran}(W), \end{array}$$

then

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

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

Remarks:

using Galerkin/one-sided projection 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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then

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

Remarks:

k = 1, standard Krylov subspace(s) of dimension $K \rightsquigarrow$ moment-matching methods/Padé approximation,

$$\frac{d^i}{ds^i}G(s_1)=\frac{d^i}{ds^i}\hat{G}(s_1), \quad i=0,\ldots, K-1(+K).$$
Intownal	atom Mada	Deduction		
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Interpolatory Model Reduction Short Introduction

Theorem (simplified) [GRIMME '97, VILLEMAGNE/SKELTON '87]

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

Remarks:

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

- dual rational Arnoldi/Lanczos [GRIMME '97],
- Iterative Rational Krylov-Algo. [ANTOULAS/BEATTIE/GUGERCIN '07].

Best \mathcal{H}_2 -norm approximation problem

Find
$$\arg \min_{\hat{G} \in \mathcal{H}_2 \text{ of order } \leq r} \|G - \hat{G}\|_2.$$

Best
$$\mathcal{H}_2$$
-norm approximation problem

Find
$$\arg\min_{\hat{G}\in\mathcal{H}_2 \text{ of order } \leq r} \|G-\hat{G}\|_2.$$

 \rightsquigarrow First-order necessary $\mathcal{H}_2\text{-}optimality$ conditions:

For SISO systems

$$G(-\mu_i) = \hat{G}(-\mu_i),$$

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

where μ_i are the poles of the reduced transfer function \hat{G} .

Best
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For MIMO systems

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

where $T^{-1}\hat{A}T = \text{diag} \{\mu_1, \dots, \mu_r\} = \text{spectral decomposition and}$ $\tilde{B} = \hat{B}^T T^{-T}, \quad \tilde{C} = \hat{C}T.$

→ tangential interpolation conditions.

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

$$\hat{G}(s) = CV \left(sI - W^{T}AV \right)^{-1} W^{T}B,$$

where V and W are given as the rational Krylov subspaces

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

$$W = \left[(-\mu_1 I - A^T)^{-1} C^T, \dots, (-\mu_r I - A^T)^{-1} C^T \right]$$

Then

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

for i = 1, ..., r as desired. \leftrightarrow iterative algorithms (IRKA/MIRIAm) that yield \mathcal{H}_2 -optimal models.

> [Gugercin et al. '06], [Bunse-Gerstner et al. '07], [Van Dooren et al. '08]

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where V and W are given as the rational Krylov subspaces

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

$$W = \left[(-\mu_1 I - A^T)^{-1} C^T, \dots, (-\mu_r I - A^T)^{-1} C^T \right].$$

Then

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

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

 \rightsquigarrow iterative algorithms (IRKA/MIRIAm) that yield \mathcal{H}_2 -optimal models.

[Gugercin et al. '06], [Bunse-Gerstner et al. '07], [Van Dooren et al. '08]

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Algorithm 1 IRKA

Input: A stable, B, C, \hat{A} stable, \hat{B} . \hat{C} . $\delta > 0$. **Output:** A^{opt}, B^{opt}, C^{opt} 1: while $(\max_{j=1,...,r} \left\{ \frac{|\mu_j - \mu_j^{\text{old}}|}{|\mu_j|} \right\} > \delta)$ do diag $\{\mu_1, \ldots, \mu_r\} := T^{-1} \hat{A} T$ = spectral decomposition, 2: $\tilde{B} = \hat{B}^* T^{-*}$ $\tilde{C} = \hat{C} T$ 3: $V = \left[(-\mu_1 I - A)^{-1} B \tilde{B}_1, \dots, (-\mu_r I - A)^{-1} B \tilde{B}_r \right]$ $W = \left[(-\mu_1 I - A^T)^{-1} C^T \tilde{C}_1, \dots, (-\mu_r I - A^T)^{-1} C^T \tilde{C}_r \right]$ 4: 5: $V = \operatorname{orth}(V), W = \operatorname{orth}(W)$ $\hat{A} = (W^*V)^{-1} W^*AV, \ \hat{B} = (W^*V)^{-1} W^*B, \ \hat{C} = CV$ 6. 7 end while 8: $A^{opt} = \hat{A}, B^{opt} = \hat{B}, C^{opt} = \hat{C}$

- Co-integration of solid fuel with silicon micromachined system.
- Goal: Ignition of solid fuel cells by electric impulse.
- Application: nano satellites.
- Thermo-dynamical model, ignition via heating an electric resistance by applying voltage source.
- Design problem: reach ignition temperature of fuel cell w/o firing neighbouring cells.
- Spatial FEM discretization of thermo-dynamical model → linear system, m = 1, p = 7.





Source: The Oberwolfach Benchmark Collection http://www.imtek.de/simulation/benchmark

Courtesy of C. Rossi, LAAS-CNRS/EU project "Micropyros".

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- axial-symmetric 2D model
- FEM discretisation using linear (quadratic) elements $\rightarrow n = 4,257$ (11,445) m = 1, p = 7.
- Reduced model computed using SPARED. modal truncation using ARPACK, and Z. Bai's PVL implementation.

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Numerical Comparison of MOR Approaches

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Topics Not Covered

- Balanced residualization (singular perturbation approximation), yields $G(0) = \hat{G}(0)$.
- Special methods for second-order (mechanical) systems.
- Extensions to bilinear and stochastic systems.
- Rational interpolation methods for nonlinear systems.
- Other MOR techniques like POD, RB.
- MOR methods for discrete-time systems.
- Extensions to descriptor systems $E\dot{x} = Ax + Bu$, E singular.
- Parametric model reduction:

$$\dot{x} = A(p)x + B(p)u, \quad y = C(p)x,$$

where $p \in \mathbb{R}^d$ is a free parameter vector; parameters should be preserved in the reduced-order model.

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