# Model Order Reduction via System Balancing 

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## Outline

Introduction

- Model Reduction for Dynamical Systems
- Application Areas
- Motivating Examples

Mathematical Basics

- Numerical Linear Algebra
- Systems and Control Theory
- Qualitative and Quantitative Study of the Approximation Error

Model Reduction by Projection

- Introduction
- Projection-based MOR Methods

Modal Truncation

- Basic Principle
- Dominant Pole Algorithm
(5) Balanced Truncation
- The basic method
- Theoretical Background
- Singular Perturbation Approximation
- Balancing-Related Methods
(6) Solving Large-Scale Matrix Equations
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- Numerical Methods for Solving Lyapunov Equations
- Solving Large-Scale Algebraic Riccati Equations
- Software


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- Model Reduction for Dynamical Systems
- Application Areas
- Motivating Examples
(2) Mathematical Basics
(3) Model Reduction by Projection
(4) Modal Truncation
(5) Balanced Truncation

6 Solving Large-Scale Matrix Equations
(7) Final Remarks

## Dynamical Systems

$$
\Sigma:\left\{\begin{array}{l}
\dot{x}(t)=f(t, x(t), u(t)), \quad x\left(t_{0}\right)=x_{0} \\
y(t)=g(t, x(t), u(t))
\end{array}\right.
$$

with

- states $x(t) \in \mathbb{R}^{n}$,
- inputs $u(t) \in \mathbb{R}^{m}$,
- outputs $y(t) \in \mathbb{R}^{q}$.



## Model Reduction for Dynamical Systems

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## Goal:

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

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## Goal:

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

## Model Reduction for Dynamical Systems

## Linear Systems

## Linear, Time-Invariant (LTI) Systems

$$
\begin{array}{lll}
\dot{x}=f(t, x, u)=A x+B u, \quad A \in \mathbb{R}^{n \times n}, & & B \in \mathbb{R}^{n \times m}, \\
y=g(t, x, u) & =C x+D u, & C \in \mathbb{R}^{q \times n},
\end{array}
$$

## Application Areas



- Resolving complex 3D geometries $\Rightarrow$ millions of degrees of freedom.
- Analysis of elastic deformations requires many simulation runs for varying external forces, in particular if the model is used in an (elastic) multi-body simulation ((E)MBS).

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


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## Application Areas

## (Optimal) Control

## Feedback Controllers

A feedback controller (dynamic compensator) is a linear system of order $N$, where

- input $=$ output of plant,
- output $=$ input of plant.

Modern (LQG-/ $\mathcal{H}_{2^{-}} / \mathcal{H}_{\infty^{-}}$) control
 design: $N \geq n$.

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

- real-time constraints,
- increasing fragility for larger $N$.
$\Longrightarrow$ reduce order of plant $(n)$ and/or controller $(N)$.
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## Application Areas

## Progressive miniaturization

- Verification of VLSI/ULSI chip design requires high number of simulations for different input signals.
- Moore's Law (1965/75) states that the number of on-chip transistors doubles each 24 months.


[^0]
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## Application Areas

## Micro Electronics/Circuit Simulation

## Progressive miniaturization

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- Increase in packing density and multilayer technology requires modeling of interconncet to ensure that thermic/electro-magnetic effects do not disturb signal transmission.

| Intel 4004 (1971) | Intel Core 2 Extreme (quad-core) (2007) |
| :--- | :--- |
| 1 layer, $10 \mu$ technology | 9 layers, 45 nm technology |
| 2,300 transistors | $>8,200,000$ transistors |
| 64 kHz clock speed | $>3 \mathrm{GHz}$ clock speed. |

## Progressive miniaturization

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Source: http://en.wikipedia.org/wiki/Image:Silicon_chip_3d.png.

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- Here: mostly MOR for linear systems, they occur in micro electronics through modified nodal analysis (MNA) for RLC networks. e.g., when
- decoupling large linear subcircuits,
- modeling transmission lines,
- modeling pin packages in VLSI chips,
- modeling circuit elements described by Maxwell's equation using partial element equivalent circuits (PEEC).


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$\rightsquigarrow$ Clear need for model reduction techniques in order to facilitate or even enable circuit simulation for current and future VLSI design.


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Standard MOR techniques in circuit simulation: Krylov subspace / Padé approximation / rational interpolation methods.

## Application Areas

Many other disciplines in computational sciences and engineering like

- computational fluid dynamics (CFD),
- computational electromagnetics,
- chemical process engineering,
- design of MEMS/NEMS (micro/nano-electrical-mechanical systems),
- computational acoustics,
- Current trend: more and more multi-physics problems, i.e., coupling of several field equations, e.g.,
- electro-thermal (e.g., bondwire heating in chip design),
- fluid-structure-interaction,


## Peter Benner and Lihong Feng.

Model Order Reduction for Coupled Problems
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## Motivating Examples

Electro-Thermic Simulation of Integrated Circuit (IC)

- Simplorer ${ }^{\circledR}$ test circuit with 2 transistors.

- Conservative thermic sub-system in Simplorer: voltage $\rightsquigarrow$ temperature, current $\rightsquigarrow$ heat flow.
- Original model: $n=270.593, m=q=2 \Rightarrow$

Computing time (on Intel Xeon dualcore 3GHz, 1 Thread):

- Main computational cost for set-up data $\approx 22 \mathrm{~min}$.
- Computation of reduced models from set-up data: 44-49sec. ( $r=20-70$ ).
- Bode plot (MATLAB on Intel Core i7, 2,67GHz, 12GB): 7.5 h for original system,$<1 \mathrm{~min}$ for reduced system.


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## Bode Plot (Amplitude)



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## Absolute Error



## Relative Error



## Motivating Examples

## A Nonlinear Model from Computational Neurosciences: the FitzHugh-Nagumo System

- Simple model for neuron (de-)activation
[Chaturantabut/Sorensen 2009]

$$
\begin{aligned}
\epsilon v_{t}(x, t) & =\epsilon^{2} v_{x x}(x, t)+f(v(x, t))-w(x, t)+g \\
w_{t}(x, t) & =h v(x, t)-\gamma w(x, t)+g
\end{aligned}
$$

with $f(v)=v(v-0.1)(1-v)$ and initial and boundary conditions

$$
\begin{array}{llr}
v(x, 0)=0, & w(x, 0)=0, & x \in[0,1] \\
v_{x}(0, t)=-i_{0}(t), & v_{x}(1, t)=0, & t \geq 0,
\end{array}
$$

where $\epsilon=0.015, h=0.5, \gamma=2, g=0.05, i_{0}(t)=50000 t^{3} \exp (-15 t)$.


Source: http://en.wikipedia.org/wiki/Neuron

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where $\epsilon=0.015, h=0.5, \gamma=2, g=0.05, i_{0}(t)=50000 t^{3} \exp (-15 t)$.

- Parameter $g$ handled as an additional input.
- Original state dimension $n=2 \cdot 400$, QBDAE dimension $N=3 \cdot 400$, reduced QBDAE dimension $r=26$, chosen expansion point $\sigma=1$.


## Motivating Examples

A Nonlinear Model from Computational Neurosciences: the FitzHugh-Nagumo System

Phase Space Diagram, $n=2 \cdot 400, r=26$


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## Numerical Linear Algebra <br> 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 $x_{i j}$ contains color information of pixel $(i, j)$.
- Memory (in single precision): $4 \cdot n_{x} \cdot n_{y}$ bytes.

```
Theorem (Schmidt-Mirsky/Eckart-Young)
```



```
\widehat{x}=\mp@subsup{\sum}{j=1}{r}\mp@subsup{\sigma}{j}{}\mp@subsup{u}{j}{}\mp@subsup{v}{j}{\top},
where }X=U\Sigma\mp@subsup{V}{}{\top}\mathrm{ is the singular value decomposition (SVD) of X.
```



Idea for dimension reduction

```
Instead of }X\mathrm{ save }\mp@subsup{u}{1}{
ur},\mp@subsup{\sigma}{1}{}\mp@subsup{v}{1}{},\ldots,\mp@subsup{\sigma}{r}{}\mp@subsup{v}{r}{
```

$\leadsto$ memory $=4 r \times\left(n_{x}+n_{y}\right)$ bytes.

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## 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}
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where $X=U \Sigma V^{\top}$ is the singular value decomposition (SVD) of $X$. The approximation error is $\|X-\widehat{X}\|_{2}=\sigma_{r+1}$.

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## Example: Image Compression by Truncated SVD

## Example: Clown



$$
\begin{gathered}
320 \times 200 \text { pixel } \\
\rightsquigarrow \quad \approx 256 \mathrm{kB}
\end{gathered}
$$

## Example: Image Compression by Truncated SVD

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- rank $r=50, \approx 104 \mathrm{kB}$


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## Dimension Reduction via SVD

## Example: Gatlinburg

Organizing committee
Gatlinburg/Householder Meeting 1964:
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$640 \times 480$ pixel, $\approx 1229$ kB

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$640 \times 480$ pixel, $\approx 1229 \mathrm{kB}$

- rank $r=100, \approx 448 \mathrm{kB}$

- rank $r=50, \approx 224 \mathrm{kB}$



## Background: Singular Value Decay

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

## Singular Values of the Image Data Matrices




## A different viewpoint

## Linear Mapping

A matrix $A \in \mathbb{R}^{\ell \times k}$ represents a linear mapping

$$
\mathcal{A}: \mathbb{R}^{k} \rightarrow \mathbb{R}^{\ell}: x \rightarrow y:=A x .
$$

The truncated SVD ignores small singular values and thus the related left and right singular vectors.

## Consequence:

- Vectors (almost) in the kernel of $A$ do not contribute to range $(A)$ and can hardly or not at all be reconstructed from the input-output relation (" $A^{-1 "}$ ) $\rightsquigarrow$ "unobservable" states.
- Vectors (almost) in range $(A)^{\perp}$ cannot be "reached" from any $x \in \mathbb{R}^{k} \rightsquigarrow$ "unreachable/uncontrollable" states.
- Hence, the truncated SVD ignores states hard to reconstruct and hard to reach.


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## Systems and Control Theory

## Definition

The Laplace transform of a time domain function $f \in L_{1, \text { loc }}$ with $\operatorname{dom}(f)=\mathbb{R}_{0}^{+}$is

$$
\mathcal{L}: f \mapsto F, \quad F(s):=\mathcal{L}\{f(t)\}(s):=\int_{0}^{\infty} e^{-s t} f(t) d t, \quad s \in \mathbb{C} .
$$

$F$ is a function in the (Laplace or) frequency domain.
Note: for frequency domain evaluations ("frequency response analysis"), one takes re $s=0$ and $\mathrm{im} s \geq 0$. Then $\omega:=\mathrm{im} s$ takes the role of a frequency (in [rad/s], i.e., $\omega=2 \pi v$ with $v$ measured in $[\mathrm{Hz}]$ ).

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Note: for ease of notation, in the following we will use lower-case letters for both, a function and its Laplace transform!

## Systems and Control Theory

The Model Reduction Problem as Approximation Problem in Frequency Domain

## Linear Systems in Frequency Domain

Application of Laplace transform $\quad(x(t) \mapsto x(s), \dot{x}(t) \mapsto s x(s))$ to linear system

$$
\dot{x}(t)=A x(t)+B u(t), \quad y(t)=C x(t)+D u(t)
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## Systems and Control Theory

The Model Reduction Problem as Approximation Problem in Frequency Domain

## Linear Systems in Frequency Domain

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Goal: Fast evaluation of mapping $u \rightarrow y$.

## Systems and Control Theory

The Model Reduction Problem as Approximation Problem in Frequency Domain

## Formulating model reduction in frequency domain

Approximate the dynamical system

$$
\begin{array}{ll}
\dot{x}=A x+B u, & A \in \mathbb{R}^{n \times n}, B \in \mathbb{R}^{n \times m} \\
y=C x+D u, & C \in \mathbb{R}^{q \times n}, D \in \mathbb{R}^{q \times m}
\end{array}
$$

by reduced-order system

$$
\begin{aligned}
& \dot{\hat{x}}=\hat{A} \hat{x}+\hat{B} u, \quad \hat{A} \in \mathbb{R}^{r \times r}, \quad \hat{B} \in \mathbb{R}^{r \times m}, \\
& \hat{y}=\hat{C} \hat{x}+\hat{D} u, \quad \hat{C} \in \mathbb{R}^{q \times r}, \quad \hat{D} \in \mathbb{R}^{q \times m}
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of order $r \ll n$, such that

$$
\|y-\hat{y}\|=\|G u-\hat{G} u\| \leq\|G-\hat{G}\| \cdot\|u\|<\text { tolerance } \cdot\|u\| .
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$\Longrightarrow$ Approximation problem:

$$
\min _{\operatorname{order}(\hat{G}) \leq r}\|G-\hat{G}\| .
$$

## Definition

A linear system

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

is stable if its transfer function $G(s)$ has all its poles in the left half plane and it is asymptotically (or Lyapunov or exponentially) stable if all poles are in the open left half plane $\mathbb{C}^{-}:=\{z \in \mathbb{C} \mid \Re(z)<0\}$.

## Lemma

Sufficient for asymptotic stability is that $A$ is asymptotically stable (or Hurwitz), i.e., the spectrum of $A$, denoted by $\Lambda(A)$, satisfies $\Lambda(A) \subset \mathbb{C}^{-}$

Note that by abuse of notation, often stable system is used for asymptotically stable systems.

## Systems and Control Theory

 Properties of linear systems
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## Systems and Control Theory

Properties of linear systems

## Questions:

- For fixed $x_{0} \in \mathbb{R}^{n}$ and some $x^{1} \in \mathbb{R}^{n}$, is there a feasible control function $u \in \mathcal{U}_{a d}$ and time $t_{1}>t_{0}=0$ such that $x\left(t_{1} ; u\right)=x^{1}$ ?
What is the set of targets $x^{1}$ reachable from $x^{0}$ ?
- For fixed $x_{1} \in \mathbb{R}^{n}$ and some $x^{0} \in \mathbb{R}^{n}$, is there a feasible control function $u \in \mathcal{U}_{a d}$ and time $t_{1}>t_{0}=0$ such that $x\left(t_{1} ; u\right)=x^{1}$ ? What is the set of initial conditions $x^{0}$ controllable to $x^{1}$ ?
E.g., $\mathcal{U}_{a d} \in\left\{C^{k}[0, T], L_{2}(0, T)\right\}$, possibly with constraints $\underline{u}(t) \leq u(t) \leq \bar{u}(t)$.


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Note: for LTI systems $\dot{x}=A x+B u$, both concepts are equivalent! E.g., $\mathcal{U}_{a d} \in\left\{C^{k}[0, T], L_{2}(0, T)\right\}$, possibly with constraints $\underline{u}(t) \leq u(t) \leq \bar{u}(t)$.

## Systems and Control Theory

 Properties of linear systems
## Definition (Controllability)

Consider the target (the state to be reached) $x^{1} \in \mathbb{R}^{n}$.
a) An LTI system with initial value $x(0)=x^{0}$ is controllable to $x^{1}$ in time $t_{1}>0$ if there exists $u \in \mathcal{U}_{a d}$ such that $x\left(t_{1} ; u\right)=x^{1}$.
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c) If the system is controllable to $x^{1}$ for all $x^{0} \in \mathbb{R}^{n}$, it is (completely) controllable.

The controllability set w.r.t. $x^{1}$ is defined as $\mathcal{C}:=\bigcup_{t_{1}>0} \mathcal{C}\left(t_{1}\right)$ where

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\mathcal{C}\left(t_{1}\right):=\left\{x^{0} \in \mathbb{R}^{n} ; \exists u \in \mathcal{U}_{a d}: x\left(t_{1} ; u\right)=x^{1}\right\} .
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In short: an LTI system is controllable $\Longleftrightarrow \mathcal{C}=\mathbb{R}^{n}$
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## Systems and Control Theory <br> Properties of linear systems

Now: characterize controllability.

## Systems and Control Theory

Properties of linear systems
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Variation of constants $\Longrightarrow$

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x(t)=e^{A t} x^{0}+\int_{0}^{t} e^{A(t-s)} B u(s) d s=e^{A t}\left(x^{0}+\int_{0}^{t} e^{-A s} B u(s) d s\right) .
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Hence, if $x^{0}$ is controllable to $x^{1}$ :

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## Systems and Control Theory

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Hence, an LTI system is controllable iff this linear system is solvable for $c \in \mathbb{R}^{n}$, i.e., iff $P\left(0, t_{1}\right)$ is invertible. (Note: $P\left(0, t_{1}\right)=P\left(0, t_{1}\right)^{T} \geq 0$ by definition!)

## Systems and Control Theory

 Properties of linear systems
## Now: characterize controllability.

## Theorem

For an LTI system defined by $(A, B) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{n \times m}$, T.F.A.E.:
a) The LTI system $\dot{x}=A x+B u$ is controllable.
b) The finite time Gramian $P\left(0, t_{1}\right)$ is spd $\forall t_{1}>0$.
c) The controllability matrix

$$
K(A, B):=\left[B, A B, A^{2} B, \ldots, A^{n-1} B\right] \in \mathbb{R}^{n \times n \cdot m}
$$

has full rank $n$. (Note: range $(K(A, B))=\mathcal{C}\left(t_{1}\right) \forall t_{1}>0$ !)
d) If $z$ is a left eigenvector of $A$, then $z^{*} B \neq 0$.
e) (Hautus test) $\operatorname{rank}([\lambda I-A, B])=n \forall \lambda \in \mathbb{C}$.

## Systems and Control Theory

Properties of linear systems
The Gramian characterization of controllability for stable systems can be based on positive definiteness of the (infinite) controllability Gramian

$$
P:=\int_{0}^{\infty} e^{A s} B B^{T} e^{A^{T} s} d s
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using congruence of $P\left(0, t_{1}\right)$ to $\int_{0}^{t_{1}} e^{A_{s}} B B^{T} e^{A^{T} s} d s$ and taking the limit $t_{1} \rightarrow \infty$.

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## Theorem

For a stable LTI system defined by $(A, B) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{n \times m}$, T.F.A.E.:
a) The LTI system $\dot{x}=A x+B u$ is controllable.
b) The controllability Gramian $P$ is positive definite.

## Systems and Control Theory

Properties of linear systems
New question: suppose we have

$$
y(t)=\tilde{y}(t)
$$

corresponding to two trajectories $x, \tilde{x}$ obtained by the same input function $u(t)$. Can we conclude that $x(0)=\tilde{x}(0)$, or even stronger, that $x(t)=\tilde{x}(t)$ for $t \leq 0, t \geq 0$ (past/future)?
(Note that $x\left(t_{0}\right)=\tilde{x}\left(t_{0}\right)$ is sufficient as trajectory uniquely determined. In other words, is the mapping $x^{0} \rightarrow y(t)$ injective?)

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## Definition (Observability)

An LTI system is reconstructable (observable) if for solution trajectories $x(t), \tilde{x}(t)$ obtained with the same input function $u$, we have

$$
\begin{aligned}
y(t) & =\tilde{y}(t) \quad \forall t \leq 0 \quad(\forall t \geq 0) \\
\Longrightarrow \quad x(t) & =\tilde{x}(t) \quad \forall t \leq 0 \quad(\forall t \geq 0)
\end{aligned}
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## Systems and Control Theory <br> Properties of linear systems

Characterization of observability/reconstructability:

## Theorem (Duality)

An LTI system is reconstructable if and only if the dual system $\dot{x}(t)=-A^{T} x(t)-C^{T} u(t)$ is controllable.

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a) The LTI system is reconstructable.
b) The LTI system is observable.
c) The observability matrix

$$
\mathcal{O}(A, C)=\left[C^{T}, A^{T} C^{T},\left(A^{2}\right)^{T} C, \ldots,\left(A^{n-1}\right)^{T} C^{T}\right]^{T} \in \mathbb{R}^{n p \times n} \text { has rank } n
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d) If $A x=\lambda x$, then $C^{\top} x \neq 0$.
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## Systems and Control Theory

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## Theorem

A stable LTI system is observable if and only if the observability Gramian

$$
Q:=\int_{0}^{\infty} e^{A^{T} t} C^{T} C e^{A t} d t
$$

is symmetric positive definite.

## Systems and Control Theory

Properties of linear systems

- Controllability/observability are sometimes too strong.
- Weaker requirement: is there $u \in \mathcal{U}_{a d}$ to steer $x_{0}$ to vicinity of $x^{1}$ ?
- For LTI systems, it suffices to consider $x^{1}=0$ !
- Hence, is there $u \in \mathcal{U}_{a d}$ so that $\lim _{t \rightarrow \infty} x(t ; u)=0\left(\forall x^{0} \in \mathbb{R}^{n}\right)$ ?
- If the answer is yes, then the LTI system is called stabilizable


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- Hence, is there $u \in \mathcal{U}_{a d}$ so that $\lim _{t \rightarrow \infty} x(t ; u)=0\left(\forall x^{0} \in \mathbb{R}^{n}\right)$ ?
- If the answer is yes, then the LTI system is called stabilizable


## Systems and Control Theory

Properties of linear systems

- Controllability/observability are sometimes too strong.
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## Theorem

For an LTI system defined by $(A, B) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{n \times m}$, T.F.A.E.:
a) The LTI system is stabilizable.
b) $\exists F \in \mathbb{R}^{m \times n}$ with $\Lambda(A+B F) \subset \mathbb{C}^{-}$.
c) If $p^{*} A=\tilde{\lambda} p^{*}$ and $\operatorname{Re}(\lambda) \geq 0$, then $p^{*} B \neq 0$.
d) $\operatorname{rank}([A-\lambda I, B])=n \quad \forall \lambda \in \mathbb{C}$ with $\operatorname{Re}(\lambda) \geq 0$.
e) In the (controllability) Kalman decomposition of $(A, B)$,

$$
V^{T} A V=\left[\begin{array}{cc}
A_{1} & A_{2} \\
0 & A_{3}
\end{array}\right], V^{T} B=\left[\begin{array}{c}
B_{1} \\
0
\end{array}\right]
$$

we have $\Lambda\left(A_{3}\right) \subset \mathbb{C}^{-}$.

## Systems and Control Theory Properties of linear systems

$\exists$ dual concept of stabilizability, analogous to duality of controllability and observability.

## Definition (Detectability)

An LTI system is detectable if for any solution $x(t)$ of $\dot{x}=A x$ with $C x(t) \equiv 0$ we have $\lim _{t \rightarrow \infty} x(t)=0$. (We can not observe all of $x$, but the unobservable part is stable.)

## Systems and Control Theory

Properties of linear systems
$\exists$ dual concept of stabilizability, analogous to duality of controllability and observability.

## Theorem

For an LTI system defined by $(A, C) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{q \times n}$, T.F.A.E.:
a) The LTI system is detectable.
b) $\left(A^{T}, C^{T}\right)$ is stabilizable.
c) $A x=\lambda x, \operatorname{Re}(\lambda) \geq 0 \Rightarrow C^{T} x \neq 0$.
d) $\operatorname{rank}\left[\begin{array}{c}\lambda I-A \\ C\end{array}\right]=n$ for all $\lambda, \operatorname{Re}(\lambda) \geq 0$.
e) In the observability Kalman decomposition of $\left(A^{T}, C^{\top}\right)$,

$$
W^{T} A W=\left[\begin{array}{cc}
A_{1} & 0 \\
A_{2} & A_{3}
\end{array}\right], C W=\left[\begin{array}{ll}
C_{1} & 0
\end{array}\right]
$$

we have $\Lambda\left(A_{3}\right) \subset \mathbb{C}^{-}$.

## Systems and Control Theory

## Realizations of Linear Systems

## Definition

For a linear (time-invariant) system

$$
\Sigma:\left\{\begin{aligned}
\dot{x}(t)=A x(t)+B u(t), & \text { with transfer function } \\
y(t)=C x(t)+D u(t), & G(s)=C(s l-A)^{-1} B+D,
\end{aligned}\right.
$$

the quadruple $(A, B, C, D) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{n \times m} \times \mathbb{R}^{q \times n} \times \mathbb{R}^{q \times m}$ is called a realization of $\Sigma$.

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## Realizations are not unique!

Transfer function is invariant under state-space transformations,

$$
\mathcal{T}:\left\{\begin{array}{ccc}
x & \rightarrow & T x, \\
(A, B, C, D) & \rightarrow & \left(T A T^{-1}, T B, C T^{-1}, D\right),
\end{array}\right.
$$

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## Realizations are not unique!

Transfer function is invariant under addition of uncontrollable/unobservable states:

$$
\begin{aligned}
\frac{d}{d t}\left[\begin{array}{c}
x \\
x_{1}
\end{array}\right] & =\left[\begin{array}{cc}
A & 0 \\
0 & A_{1}
\end{array}\right]\left[\begin{array}{c}
x \\
x_{1}
\end{array}\right]+\left[\begin{array}{c}
B \\
B_{1}
\end{array}\right] u(t), \quad y(t)=\left[\begin{array}{ll}
C & 0
\end{array}\right]\left[\begin{array}{c}
x \\
x_{1}
\end{array}\right]+D u(t), \\
\frac{d}{d t}\left[\begin{array}{c}
x \\
x_{2}
\end{array}\right] & =\left[\begin{array}{cc}
A & 0 \\
0 & A_{2}
\end{array}\right]\left[\begin{array}{c}
x \\
x_{2}
\end{array}\right]+\left[\begin{array}{c}
B \\
0
\end{array}\right] u(t), \quad y(t)=\left[\begin{array}{ll}
C & C_{2}
\end{array}\right]\left[\begin{array}{c}
x \\
x_{2}
\end{array}\right]+D u(t),
\end{aligned}
$$

for arbitrary $A_{j} \in \mathbb{R}^{n_{j} \times n_{j}}, j=1,2, B_{1} \in \mathbb{R}^{n_{1} \times m}, C_{2} \in \mathbb{R}^{q \times n_{2}}$ and any $n_{1}, n_{2} \in \mathbb{N}$.

## Systems and Control Theory

## Realizations of Linear Systems

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the quadruple $(A, B, C, D) \in \mathbb{R}^{n \times n} \times \mathbb{R}^{n \times m} \times \mathbb{R}^{q \times n} \times \mathbb{R}^{q \times m}$ is called a realization of $\Sigma$.

## Realizations are not unique!

Hence,

$$
\begin{array}{ll}
(A, B, C, D), & \left(\left[\begin{array}{cc}
A & 0 \\
0 & A_{1}
\end{array}\right],\left[\begin{array}{c}
B \\
B_{1}
\end{array}\right],\left[\begin{array}{ll}
C & 0
\end{array}\right], D\right), \\
\left(T A T^{-1}, T B, C T^{-1}, D\right), & \left(\left[\begin{array}{cc}
A & 0 \\
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\end{array}
$$

are all realizations of $\Sigma$ !

## Systems and Control Theory

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## Definition

The McMillan degree of $\Sigma$ is the unique minimal number $\hat{n} \geq 0$ of states necessary to describe the input-output behavior completely. A minimal realization is a realization $(\hat{A}, \hat{B}, \hat{C}, \hat{D})$ of $\Sigma$ with order $\hat{n}$.

## Systems and Control Theory

## Realizations of Linear Systems

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## Definition

The McMillan degree of $\Sigma$ is the unique minimal number $\hat{n} \geq 0$ of states necessary to describe the input-output behavior completely. A minimal realization is a realization $(\hat{A}, \hat{B}, \hat{C}, \hat{D})$ of $\Sigma$ with order $\hat{n}$.

## Theorem

A realization $(A, B, C, D)$ of a linear system is minimal $\Longleftrightarrow$ $(A, B)$ is controllable and $(A, C)$ is observable.

## Systems and Control Theory

## Balanced Realizations

## Definition

A realization $(A, B, C, D)$ of a linear system $\Sigma$ is balanced if its infinite controllability/observability Gramians $P / Q$ satisfy

$$
P=Q=\operatorname{diag}\left\{\sigma_{1}, \ldots, \sigma_{n}\right\} \quad \text { (w.l.o.g. } \sigma_{j} \geq \sigma_{j+1}, j=1, \ldots, n-1 \text { ). }
$$

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When does a balanced realization exist?

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$$

When does a balanced realization exist?
Assume $A$ to be Hurwitz, i.e. $\Lambda(A) \subset \mathbb{C}^{-}$. Then:

## Theorem

Given a stable minimal linear system $\Sigma:(A, B, C, D)$, a balanced realization is obtained by the state-space transformation with

$$
T_{b}:=\Sigma^{-\frac{1}{2}} V^{\top} R,
$$

where $P=S^{T} S, Q=R^{T} R$ (e.g., Cholesky decompositions) and $S R^{T}=U \Sigma V^{T}$ is the SVD of $S R^{T}$.

Proof. Easy.

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$$

$\sigma_{1}, \ldots, \sigma_{n}$ are the Hankel singular values of $\Sigma$.
Note: $\sigma_{1}, \ldots, \sigma_{n} \geq 0$ as $P, Q \geq 0$ by definition, and $\sigma_{1}, \ldots, \sigma_{n}>0$ in case of minimality!

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## Theorem

The infinite controllability/observability Gramians $P / Q$ satisfy the Lyapunov equations

$$
A P+P A^{T}+B B^{T}=0, \quad A^{T} Q+Q A+C^{T} C=0 .
$$

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Proof. Exercise!

## Definition

A realization ( $A, B, C, D$ ) of a stable linear system $\Sigma$ is balanced if its infinite controllability/observability Gramians $P / Q$ satisfy

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Note: $\sigma_{1}, \ldots, \sigma_{n} \geq 0$ as $P, Q \geq 0$ by definition, and $\sigma_{1}, \ldots, \sigma_{n}>0$ in case of minimality!

## Theorem

The Hankel singular values (HSVs) of a stable minimal linear system are system invariants, i.e. they are unaltered by state-space transformations!

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Proof. In balanced coordinates, the HSVs are $\Lambda(P Q)^{\frac{1}{2}}$. Now let

$$
(\hat{A}, \hat{B}, \hat{C}, D)=\left(T A T^{-1}, T B, C T^{-1}, D\right)
$$

be any transformed realization with associated controllability Lyapunov equation

$$
0=\hat{A} \hat{P}+\hat{P} \hat{A}^{T}+\hat{B} \hat{B}^{T}=T A T^{-1} \hat{P}+\hat{P} T^{-T} A^{T} T^{T}+T B B^{T} T^{T} .
$$

This is equivalent to

$$
0=A\left(T^{-1} \hat{P} T^{-T}\right)+\left(T^{-1} \hat{P} T^{-T}\right) A^{T}+B B^{T}
$$

The uniqueness of the solution of the Lyapunov equation implies that $\hat{P}=T P T^{T}$ and, analogously, $\hat{Q}=T^{-T} Q T^{-1}$. Therefore,

$$
\hat{P} \hat{Q}=T P Q T^{-1}
$$

showing that $\Lambda(\hat{P} \hat{Q})=\Lambda(P Q)=\left\{\sigma_{1}^{2}, \ldots, \sigma_{n}^{2}\right\}$.

## Systems and Control Theory Balanced Realizations

## Definition

A realization $(A, B, C, D)$ of a stable linear system $\Sigma$ is balanced if its infinite controllability/observability Gramians $P / Q$ satisfy

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\left.P=Q=\operatorname{diag}\left\{\sigma_{1}, \ldots, \sigma_{n}\right\} \quad \text { (w.l.o.g. } \sigma_{j} \geq \sigma_{j+1}, j=1, \ldots, n-1\right)
$$

$\sigma_{1}, \ldots, \sigma_{n}$ are the Hankel singular values of $\Sigma$.
Note: $\sigma_{1}, \ldots, \sigma_{n} \geq 0$ as $P, Q \geq 0$ by definition, and $\sigma_{1}, \ldots, \sigma_{n}>0$ in case of minimality!

## Remark

For non-minimal systems, the Gramians can also be transformed into diagonal matrices with the leading $\hat{n} \times \hat{n}$ submatrices equal to $\operatorname{diag}\left(\sigma_{1}, \ldots, \sigma_{\hat{n}}\right)$, and

$$
\hat{P} \hat{Q}=\operatorname{diag}\left(\sigma_{1}^{2}, \ldots, \sigma_{\hat{n}}^{2}, 0, \ldots, 0\right)
$$

see [Laub/Heath/Paige/Ward 1987, Tombs/Postlethwaite 1987].

## Qualitative and Quantitative Study of the Approximation Error

 System NormsConsider transfer function

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

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

$$
\|u\|_{2}^{2}:=\frac{1}{2 \pi} \int_{-\infty}^{\infty} u(\jmath \omega)^{H} u(\jmath \omega) d \omega .
$$

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

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Assume $A$ (asympotically) stable: $\Lambda(A) \subset \mathbb{C}^{-}:=\{z \in \mathbb{C}:$ re $z<0\}$. Then for all $s \in \mathbb{C}^{+} \cup \jmath \mathbb{R},\|G(s)\| \leq M<\infty \Rightarrow$

$$
\int_{-\infty}^{\infty} y(\jmath \omega)^{H} y(\jmath \omega) d \omega=\int_{-\infty}^{\infty} u(\jmath \omega)^{H} G(\jmath \omega)^{H} G(\jmath \omega) u(\jmath \omega) d \omega
$$

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

## Qualitative and Quantitative Study of the Approximation Error System Norms

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$$
\begin{aligned}
\int_{-\infty}^{\infty} y(\jmath \omega)^{H} y(\jmath \omega) d \omega & =\int_{-\infty}^{\infty} u(\jmath \omega)^{H} G(\jmath \omega)^{H} G(\jmath \omega) u(\jmath \omega) d \omega \\
& =\int_{-\infty}^{\infty}\|G(\jmath \omega) u(\jmath \omega)\|^{2} d \omega \leq \int_{-\infty}^{\infty} M^{2}\|u(\jmath \omega)\|^{2} d \omega
\end{aligned}
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& =\int_{-\infty}^{\infty}\|G(\jmath \omega) u(\jmath \omega)\|^{2} d \omega \leq \int_{-\infty}^{\infty} M^{2}\|u(\jmath \omega)\|^{2} d \omega \\
& =M^{2} \int_{-\infty}^{\infty} u(\jmath \omega)^{H} u(\jmath \omega) d \omega<\infty .
\end{aligned}
$$

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

## Qualitative and Quantitative Study of the Approximation Error

## System Norms

Consider transfer function

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& =M^{2} \int_{-\infty}^{\infty} u(\jmath \omega)^{H} u(\jmath \omega) d \omega<\infty
\end{aligned}
$$

$$
\Longrightarrow y \in \mathcal{L}_{2}^{q} \cong L_{2}^{q}(-\infty, \infty)
$$

## Qualitative and Quantitative Study of the Approximation Error

 System NormsConsider transfer function

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$$

Assume $A$ (asympotically) stable: $\Lambda(A) \subset \mathbb{C}^{-}:=\{z \in \mathbb{C}:$ re $z<0\}$. Consequently, the 2-induced operator norm

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

is well defined. It can be shown that

$$
\|G\|_{\infty}=\sup _{\omega \in \mathbb{R}}\|G(\jmath \omega)\|=\sup _{\omega \in \mathbb{R}} \sigma_{\max }(G(\jmath \omega)) .
$$

## Qualitative and Quantitative Study of the Approximation Error

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is well defined. It can be shown that

$$
\|G\|_{\infty}=\sup _{\omega \in \mathbb{R}}\|G(\jmath \omega)\|=\sup _{\omega \in \mathbb{R}} \sigma_{\max }(G(\jmath \omega))
$$

Sketch of proof:
$\|G(\jmath \omega) u(\jmath \omega)\| \leq\|G(\jmath \omega)\|\|u(\jmath \omega)\| \Rightarrow " \leq "$.
Construct $u$ with $\|G u\|_{2}=\sup _{\omega \in \mathbb{R}}\|G(\jmath \omega)\|\|u\|_{2}$.

## Qualitative and Quantitative Study of the Approximation Error

 System NormsConsider transfer function

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

## Hardy space $\mathcal{H}_{\infty}$

Function space of matrix-/scalar-valued functions that are analytic and bounded in $\mathbb{C}^{+}$.
The $\mathcal{H}_{\infty}$-norm is

$$
\|F\|_{\infty}:=\sup _{\text {res>0 }} \sigma_{\max }(F(s))=\sup _{\omega \in \mathbb{R}} \sigma_{\max }(F(\jmath \omega))
$$

Stable transfer functions are in the Hardy spaces

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

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Consequence of Parseval identity/Plancherel Theorem

$$
L_{2}(-\infty, \infty) \cong \mathcal{L}_{2}, \quad L_{2}(0, \infty) \cong \mathcal{H}_{2}
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Consequently, 2-norms in time and frequency domains coincide!

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## $\mathcal{H}_{\infty}$ approximation error

Reduced-order model $\Rightarrow$ transfer function $\hat{G}(s)=\hat{C}\left(s I_{r}-\hat{A}\right)^{-1} \hat{B}+\hat{D}$.

$$
\|y-\hat{y}\|_{2}=\|G u-\hat{G} u\|_{2} \leq\|G-\hat{G}\|_{\infty}\|u\|_{2} .
$$

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

## Qualitative and Quantitative Study of the Approximation Error

 System NormsConsider stable transfer function

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

## Hardy space $\mathcal{H}_{2}$

Function space of matrix-/scalar-valued functions that are analytic $\mathbb{C}^{+}$and bounded w.r.t. the $\mathcal{H}_{2}$-norm

$$
\begin{aligned}
\|F\|_{2} & :=\frac{1}{2 \pi}\left(\sup _{\operatorname{re\sigma } \sigma 0} \int_{-\infty}^{\infty}\|F(\sigma+\jmath \omega)\|_{F}^{2} d \omega\right)^{\frac{1}{2}} \\
& =\frac{1}{2 \pi}\left(\int_{-\infty}^{\infty}\|F(\jmath \omega)\|_{F}^{2} d \omega\right)^{\frac{1}{2}}
\end{aligned}
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\|F\|_{2}=\frac{1}{2 \pi}\left(\int_{-\infty}^{\infty}\|F(\jmath \omega)\|_{F}^{2} d \omega\right)^{\frac{1}{2}}
$$

## $\mathcal{H}_{2}$ approximation error for impulse response $\left(u(t)=u_{0} \delta(t)\right)$

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

$$
\|y-\hat{y}\|_{2}=\left\|G u_{0} \delta-\hat{G} u_{0} \delta\right\|_{2} \leq\|G-\hat{G}\|_{2}\left\|u_{0}\right\| .
$$

$\Longrightarrow$ compute reduced-order model such that $\|G-\hat{G}\|_{2}<$ tol!

## Qualitative and Quantitative Study of the Approximation Error

 System NormsConsider stable transfer function

$$
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$$
\|F\|_{2}=\frac{1}{2 \pi}\left(\int_{-\infty}^{\infty}\|F(\jmath \omega)\|_{F}^{2} d \omega\right)^{\frac{1}{2}}
$$

## Theorem (Practical Computation of the $\mathcal{H}_{2}$-norm)

$$
\|F\|_{2}^{2}=\operatorname{tr}\left(B^{\top} Q B\right)=\operatorname{tr}\left(C P C^{\top}\right),
$$

where $P, Q$ are the controllability and observability Gramians of the corresponding LTI system.

## Output errors in time-domain

$$
\begin{aligned}
\|y-\hat{y}\|_{2} & \leq\|G-\hat{G}\|_{\infty}\|u\|_{2}
\end{aligned} \quad \Longrightarrow\|G-\hat{G}\|_{\infty}<\text { tol }
$$

## Qualitative and Quantitative Study of the Approximation Error Approximation Problems

## Output errors in time-domain

$$
\begin{aligned}
\|y-\hat{y}\|_{2} & \leq\|G-\hat{G}\|_{\infty}\|u\|_{2}
\end{aligned} \quad \Longrightarrow\|G-\hat{G}\|_{\infty}<\text { tol }
$$

| $\mathcal{H}_{\infty}$-norm | best approximation problem for given reduced order $r$ in <br> general open; balanced truncation yields suboptimal solu- <br> tion with computable $\mathcal{H}_{\infty}$-norm bound. |
| :--- | :--- |
| $\mathcal{H}_{2}$-norm | necessary conditions for best approximation known; (local) <br> optimizer computable with iterative rational Krylov algo- <br> rithm (IRKA) |
| Hankel-norm <br> $\\|G\\|_{H}:=\sigma_{\max }$ | optimal Hankel norm approximation (AAK theory). |

## Qualitative and Quantitative Study of the Approximation Error <br> \section*{Computable error measures}

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

## Other measures

- absolute errors $\left\|G\left(\jmath \omega_{j}\right)-\hat{G}\left(\jmath \omega_{j}\right)\right\|_{2},\left\|G\left(\jmath \omega_{j}\right)-\hat{G}\left(\jmath \omega_{j}\right)\right\|_{\infty}\left(j=1, \ldots, N_{\omega}\right)$;
- relative errors $\frac{\left\|G\left(\jmath \omega_{j}\right)-\hat{G}\left(\jmath \omega_{j}\right)\right\|_{2}}{\left\|G\left(\jmath \omega_{j}\right)\right\|_{2}}, \frac{\left\|G\left(\jmath \omega_{j}\right)-\hat{G}\left(\jmath \omega_{j}\right)\right\|_{\infty}}{\left\|G\left(\jmath \omega_{j}\right)\right\| \infty}$;
- "eyeball norm", i.e. look at frequency response/Bode (magnitude) plot: for SISO system, log-log plot frequency vs. $|G(\jmath \omega)|$ (or $|G(\jmath \omega)-\hat{G}(\jmath \omega)|)$ in decibels, $1 \mathrm{~dB} \simeq 20 \log _{10}$ (value).
For MIMO systems, $q \times m$ array of plots $G_{i j}$.


(c) Peter Benner, MOR via System Balancing


## Outline

(1) Introduction
(2) Mathematical Basics
(3) Model Reduction by Projection

- Introduction
- Projection-based MOR Methods
(4) Modal Truncation
(5) Balanced Truncation

6 Solving Large-Scale Matrix Equations
(7) Final Remarks

## Model Reduction by Projection

## Goals

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

$$
\|y-\hat{y}\|<\text { tolerance } \cdot\|u\| \quad \forall u \in L_{2}\left(\mathbb{R}, \mathbb{R}^{m}\right)
$$

$\Longrightarrow$ Need computable error bound/estimate!

- Preserve physical properties:


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("system does not generate energy").


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$$
\int_{-\infty}^{t} u(\tau)^{T} y(\tau) d \tau \geq 0 \quad \forall t \in \mathbb{R}, \quad \forall u \in L_{2}\left(\mathbb{R}, \mathbb{R}^{m}\right)
$$

("system does not generate energy").

## Model Reduction by Projection

Projection Basics

## Definition (Projector)

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

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## Lemma (Projector Properties)

- 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., $P v=v \forall v \in \mathcal{V}$
- $I-P$ is the complementary projector onto $\operatorname{ker} P$
- If $\mathcal{V}$ is an $A$-invariant subspace corresponding to a subset of $A^{\prime}$ 's spectrum, then we call $P$ a spectral projector.
- Let $\mathcal{W} \subset \mathbb{R}^{n}$ be another $r$-dimensional subspace and $W=\left[w_{1}, \ldots, w_{r}\right]$ be a basis matrix for $\mathcal{W}$, then $P=V\left(W^{\top} V\right)^{-1} W^{\top}$ is an oblique projector onto $\mathcal{V}$ along $\mathcal{W}$.


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## Model Reduction by Projection

Projection-based MOR Methods

## Methods:

(1) Modal Truncation
(2) Balanced Truncation
(3) Rational Interpolation (Padé-Approximation and (rational) Krylov Subspace Methods)
( ( many more...
Joint feature of these methods:
computation of reduced-order model (ROM) by projection!

## Model Reduction by Projection

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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

$$
\operatorname{range}(V)=\mathcal{V}, \quad \text { range }(W)=\mathcal{W}, \quad W^{T} V=I_{r}
$$

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

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

Important observation:

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

$$
W^{T}(\dot{\tilde{x}}-A \tilde{x}-B u)=W^{T}\left(V W^{T} \dot{x}-A V W^{T} x-B u\right)
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& =\dot{\hat{x}}-\hat{A} \hat{x}-\hat{B} u=0
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## Outline

(1) Introduction
(2) Mathematical Basics
(3) Model Reduction by Projection
(4) Modal Truncation

- Basic Principle
- Dominant Pole Algorithm
(5) Balanced Truncation

6 Solving Large-Scale Matrix Equations
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## Modal Truncation

## Basic method:

Assume $A$ is diagonalizable, $T^{-1} A T=D_{A}$, project state-space onto $A$-invariant subspace $\mathcal{V}=\operatorname{span}\left(t_{1}, \ldots, t_{r}\right), t_{k}=$ eigenvectors corresp. to "dominant" modes / eigenvalues of $A$. Then with

$$
V=T(:, 1: r)=\left[t_{1}, \ldots, t_{r}\right], \quad \tilde{W}^{H}=T^{-1}(1: r,:), \quad W=\tilde{W}\left(V^{H} \tilde{W}\right)^{-1}
$$

reduced-order model is

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\hat{A}:=W^{H} A V=\operatorname{diag}\left\{\lambda_{1}, \ldots, \lambda_{r}\right\}, \quad \hat{B}:=W^{H} B, \quad \hat{C}=C V
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Also computable by truncation:

$$
T^{-1} A T=\left[\begin{array}{cc}
\hat{A} & \\
& A_{2}
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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:

$$
\|G-\hat{G}\|_{\infty} \leq\left\|C_{2}\right\|\left\|B_{2}\right\| \frac{1}{\min _{\lambda \in \Lambda\left(A_{2}\right)}|\operatorname{Re}(\lambda)|}
$$

Proof:

$$
\left.\begin{array}{rl}
G(s) & =C(s I-A)^{-1} B+D=C T T^{-1}(s I-A)^{-1} T T^{-1} B+D \\
& =C T\left(s I-T^{-1} A T\right)^{-1} T^{-1} B+D \\
& =\left[\hat{C}, C_{2}\right]\left[\begin{array}{cc}
\left(s I_{r}-\hat{A}\right)^{-1} & \\
& =G(s)+C_{2}\left(s I_{n-r}-A_{2}\right)^{-1} B_{2},
\end{array}\right]\left[\begin{array}{c}
\hat{B} \\
B_{2}
\end{array}\right]+D \\
& =\hat{G}(s)^{-1}
\end{array}\right]
$$

## Modal Truncation

## Basic method:

$$
T^{-1} A T=\left[\begin{array}{cc}
\hat{A} & \\
& A_{2}
\end{array}\right], \quad T^{-1} B=\left[\begin{array}{c}
\hat{B} \\
B_{2}
\end{array}\right], \quad C T=\left[\hat{C}, C_{2}\right], \quad \hat{D}=D .
$$

## Properties:

## Error bound:

$$
\|G-\hat{G}\|_{\infty} \leq\left\|C_{2}\right\|\left\|B_{2}\right\| \frac{1}{\min _{\lambda \in \Lambda\left(A_{2}\right)}|\operatorname{Re}(\lambda)|}
$$

Proof:

$$
G(s)=\hat{G}(s)+C_{2}\left(s I_{n-r}-A_{2}\right)^{-1} B_{2},
$$

observing that $\|G-\hat{G}\|_{\infty}=\sup _{\omega \in \mathbb{R}} \sigma_{\max }\left(C_{2}\left(\jmath \omega I_{n-r}-A_{2}\right)^{-1} B_{2}\right)$, and

$$
C_{2}\left(\jmath \omega I_{n-r}-A_{2}\right)^{-1} B_{2}=C_{2} \operatorname{diag}\left(\frac{1}{\jmath \omega-\lambda_{r+1}}, \ldots, \frac{1}{\jmath \omega-\lambda_{n}}\right) B_{2} .
$$

## Modal Truncation

## Basic method:

Assume $A$ is diagonalizable, $T^{-1} A T=D_{A}$, project state-space onto $A$-invariant subspace $\mathcal{V}=\operatorname{span}\left(t_{1}, \ldots, t_{r}\right), t_{k}=$ eigenvectors corresp. to "dominant" modes / eigenvalues of $A$. Then reduced-order model is

$$
\hat{A}:=W^{H} A V=\operatorname{diag}\left\{\lambda_{1}, \ldots, \lambda_{r}\right\}, \quad \hat{B}:=W^{H} B, \quad \hat{C}=C V
$$

Also computable by truncation:

$$
T^{-1} A T=\left[\begin{array}{cc}
\hat{A} & \\
& A_{2}
\end{array}\right], \quad T^{-1} B=\left[\begin{array}{c}
\hat{B} \\
B_{2}
\end{array}\right], \quad C T=\left[\hat{C}, C_{2}\right], \quad \hat{D}=D
$$

## 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.


## Basic Principle

## Example

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

## Bode plots of transfer functions and error function




MATLAB ${ }^{\circledR}$ demo

## Basic Principle

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## Bode plots of transfer functions and error function




MATLAB ${ }^{\circledR}$ demo.

## Basic Principle

## Extensions

## Base enrichment

Static modes are defined by setting $\dot{x}=0$ and assuming unit loads, i.e., $u(t) \equiv e_{j}, j=1, \ldots, m$ :

$$
0=A x(t)+B e_{j} \quad \Longrightarrow \quad x(t) \equiv-A^{-1} b_{j} .
$$

Projection subspace $\mathcal{V}$ is then augmented by $A^{-1}\left[b_{1}, \ldots, b_{m}\right]=A^{-1} B$. Interpolation-projection framework $\Longrightarrow G(0)=\hat{G}(0)$ !
If two sided projection is used, complimentary subspace can be augmented by $A^{-T} C^{T} \Longrightarrow G^{\prime}(0)=\hat{G}^{\prime}(0)$ !
Note: if $m \neq q$, add random vectors or delete some of the columns in $A^{-T} C^{T}$.

## Basic Principle

## Extensions

## Guyan reduction (static condensation)

Partition states in masters $x_{1} \in \mathbb{R}^{r}$ and slaves $x_{2} \in \mathbb{R}^{n-r}$ (FEM terminology) Assume stationarity, i.e., $\dot{x}=0$ and solve for $x_{2}$ in

$$
\begin{aligned}
0 & =\left[\begin{array}{ll}
A_{11} & A_{12} \\
A_{21} & A_{22}
\end{array}\right]\left[\begin{array}{l}
x_{1} \\
x_{2}
\end{array}\right]+\left[\begin{array}{c}
B_{1} \\
B_{2}
\end{array}\right] u \\
\Rightarrow \quad x_{2} & =-A_{22}^{-1} A_{21} x_{1}-A_{22}^{-1} B_{2} u .
\end{aligned}
$$

Inserting this into the first part of the dynamic system

$$
\dot{x}_{1}=A_{11} x_{1}+A_{12} x_{2}+B_{1} u, \quad y=C_{1} x_{1}+C_{2} x_{2}
$$

then yields the reduced-order model

$$
\begin{aligned}
\dot{x}_{1} & =\left(A_{11}-A_{12} A_{22}^{-1} A_{21}\right) x_{1}+\left(B_{1}-A_{12} A_{22}^{-1} B_{2}\right) u \\
y & =\left(C_{1}-C_{2} A_{22}^{-1} A_{21}\right) x_{1}-C_{2} A_{22}^{-1} B_{2} u .
\end{aligned}
$$

## Modal Truncation

## Pole-Residue Form of Transfer Function

Consider partial fraction expansion of transfer function with $D=0$ :

$$
G(s)=\sum_{k=1}^{n} \frac{R_{k}}{s-\lambda_{k}}
$$

with the residues $R_{k}:=\left(C x_{k}\right)\left(y_{k}^{H} B\right) \in \mathbb{C}^{q \times m}$.

## Modal Truncation

## Dominant Pole Algorithm

## Pole-Residue Form of Transfer Function

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$$

with the residues $R_{k}:=\left(C x_{k}\right)\left(y_{k}^{H} B\right) \in \mathbb{C}^{q \times m}$.
Note: this follows using the spectral decomposition $A=X D X^{-1}$, with $X=\left[x_{1}, \ldots, x_{n}\right]$ the right and $X^{-1}=: Y=\left[y_{1}, \ldots, y_{n}\right]^{H}$ the left eigenvector matrices:

$$
\begin{aligned}
G(s) & =C\left(s l-X D X^{-1}\right)^{-1} B=C X\left(s I-\operatorname{diag}\left\{\lambda_{1}, \ldots, \lambda_{n}\right\}\right)^{-1} Y B \\
& =\left[C x_{1}, \ldots, C_{x_{n}}\right]\left[\begin{array}{ccc}
\frac{1}{s-\lambda_{1}} & & \\
& \ddots & \\
& & \left.\frac{1}{s-\lambda_{n}}\right]\left[\begin{array}{l}
y_{1}^{H} B \\
\vdots \\
y_{n}^{H} B
\end{array}\right] \\
& =\sum_{k=1}^{n} \frac{\left(C x_{k}\right)\left(y_{k}^{H} B\right)}{s-\lambda_{k}} .
\end{array} .\right.
\end{aligned}
$$

## Modal Truncation

Dominant Pole Algorithm

## Pole-Residue Form of Transfer Function

Consider partial fraction expansion of transfer function with $D=0$ :

$$
G(s)=\sum_{k=1}^{n} \frac{R_{k}}{s-\lambda_{k}}
$$

with the residues $R_{k}:=\left(C x_{k}\right)\left(y_{k}^{H} B\right) \in \mathbb{C}^{q \times m}$.
Note: $R_{k}=\left(C x_{k}\right)\left(y_{k}^{H} B\right)$ are the residues of $G$ in the sense of the residue theorem of complex analysis:

$$
\begin{aligned}
\operatorname{res}\left(G, \lambda_{\ell}\right)=\lim _{s \rightarrow \lambda_{\ell}}\left(s-\lambda_{\ell}\right) G(s)=\sum_{k=1}^{n} & \underbrace{\lim _{s \rightarrow \lambda_{\ell}} \frac{s-\lambda_{\ell}}{s-\lambda_{k}}} \quad R_{k}=R_{\ell} . \\
= & \left\{\begin{array}{l}
0 \text { for } k \neq \ell \\
1 \text { for } k=\ell
\end{array}\right.
\end{aligned}
$$

## Modal Truncation

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with the residues $R_{k}:=\left(C x_{k}\right)\left(y_{k}^{H} B\right) \in \mathbb{C}^{q \times m}$.
As projection basis use spaces spanned by right/left eigenvectors corresponding to dominant poles, i.e.. $\left(\lambda_{j}, x_{j}, y_{j}\right)$ with largest

$$
\left\|R_{k}\right\| / \mid \text { re }\left(\lambda_{k}\right) \mid .
$$

## Modal Truncation

## Dominant Pole Algorithm

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$$

## Remark

The dominant modes have most important influence on the input-output behavior of the system and are responsible for the "peaks"' in the frequency response.

## Dominant Poles

Random SISO Example ( $B, C^{T} \in \mathbb{R}^{n}$ )


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## Dominant Poles

Random SISO Example ( $B, C^{T} \in \mathbb{R}^{n}$ )

Algorithms for computing dominant poles and eigenvectors:

- Subspace Accelerated Dominante Pole Algorithm (SADPA),
- Rayleigh-Quotient-Iteration (RQI),
- Jacobi-Davidson-Method.


## Outline

(1) Introduction
(2) Mathematical Basics
(3) Model Reduction by Projection
(4) Modal Truncation
(5) Balanced Truncation

- The basic method
- Theoretical Background
- Singular Perturbation Approximation
- Balancing-Related Methods
(6) Solving Large-Scale Matrix Equations
(7) Final Remarks


## Balanced Truncation

## Basic principle:

- Recall: a stable system $\Sigma$, realized by $(A, B, C, D)$, is called balanced, if the Gramians, i.e., solutions $P, Q$ of the Lyapunov equations

$$
A P+P A^{T}+B B^{T}=0, \quad A^{T} Q+Q A+C^{T} C=0
$$

satisfy: $P=Q=\operatorname{diag}\left(\sigma_{1}, \ldots, \sigma_{n}\right)$ with $\sigma_{1} \geq \sigma_{2} \geq \ldots \geq \sigma_{n}>0$.

- $\wedge(P Q)^{\frac{1}{2}}=\left\{\sigma_{1}, \ldots, \sigma_{n}\right\}$ are the Hankel singular values (HSVs) of $\Sigma$.


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- $\Lambda(P Q)^{\frac{1}{2}}=\left\{\sigma_{1}, \ldots, \sigma_{n}\right\}$ are the Hankel singular values (HSVs) of $\Sigma$.
- Compute balanced realization of the system via state-space transformation

$$
\begin{aligned}
\mathcal{T}:(A, B, C, D) & \mapsto\left(T A T^{-1}, T B, C T^{-1}, D\right) \\
& =\left(\left[\begin{array}{ll}
A_{11} & A_{12} \\
A_{21} & A_{22}
\end{array}\right],\left[\begin{array}{l}
B_{1} \\
B_{2}
\end{array}\right],\left[\begin{array}{ll}
C_{1} & C_{2}
\end{array}\right], D\right)
\end{aligned}
$$

- Truncation $\rightsquigarrow(\hat{A}, \hat{B}, \hat{C}, \hat{D}):=\left(A_{11}, B_{1}, C_{1}, D\right)$.


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## Balanced Truncation

## Motivation:

The HSVs $\wedge(P Q)^{\frac{1}{2}}=\left\{\sigma_{1}, \ldots, \sigma_{n}\right\}$ are system invariants: they are preserved under

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$$

in transformed coordinates, the Gramians satisfy

$$
\begin{aligned}
& \qquad \begin{aligned}
\left(T A T^{-1}\right)\left(T P T^{T}\right)+\left(T P T^{T}\right)\left(T A T^{-1}\right)^{T}+(T B)(T B)^{T}=0, \\
\left(T A T^{-1}\right)^{T}\left(T^{-T} Q T^{-1}\right)+\left(T^{-T} Q T^{-1}\right)\left(T A T^{-1}\right)+\left(C T^{-1}\right)^{T}\left(C T^{-1}\right)=0 \\
\Rightarrow\left(T P T^{T}\right)\left(T^{-T} Q T^{-1}\right)=T P Q T^{-1},
\end{aligned} \\
& \text { hence } \Lambda(P Q)=\Lambda\left(\left(T P T^{T}\right)\left(T^{-T} Q T^{-1}\right)\right) .
\end{aligned}
$$

## Balanced Truncation

## Implementation: SR Method

(1) Compute (Cholesky) factors of the Gramians, $P=S^{\top} S, Q=R^{T} R$.
(0) Compute SVD $S R^{T}=\left[U_{1}, U_{2}\right]\left[\begin{array}{ll}\Sigma_{1} & \\ & \Sigma_{2}\end{array}\right]\left[\begin{array}{l}V_{1}^{\top} \\ V_{2}^{\top}\end{array}\right]$

- ROM is $\left(W^{T} A V, W^{T} B, C V, D\right)$, where

$$
W=R^{T} V_{1} \Sigma_{1}^{-\frac{1}{2}}, \quad V=S^{\top} U_{1} \Sigma_{1}^{-\frac{1}{2}} .
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$$
W=R^{T} V_{1} \Sigma_{1}^{-\frac{1}{2}}, \quad V=S^{T} U_{1} \Sigma_{1}^{-\frac{1}{2}}
$$

Note:

$$
V^{\top} W=\left(\Sigma_{1}^{-\frac{1}{2}} U_{1}^{T} S\right)\left(R^{T} V_{1} \Sigma_{1}^{-\frac{1}{2}}\right)
$$

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Note:

$$
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Note:

$$
\begin{aligned}
V^{\top} W & =\left(\Sigma_{1}^{-\frac{1}{2}} U_{1}^{T} S\right)\left(R^{T} V_{1} \Sigma_{1}^{-\frac{1}{2}}\right)=\Sigma_{1}^{-\frac{1}{2}} U_{1}^{T} U \Sigma V^{\top} V_{1} \Sigma_{1}^{-\frac{1}{2}} \\
& =\Sigma_{1}^{-\frac{1}{2}}\left[I_{r}, 0\right]\left[\begin{array}{cc}
\Sigma_{1} & \\
& \Sigma_{2}
\end{array}\right]\left[\begin{array}{c}
I_{r} \\
0
\end{array}\right] \Sigma_{1}^{-\frac{1}{2}}
\end{aligned}
$$

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$$
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\Sigma_{1} & \\
& \Sigma_{2}
\end{array}\right]\left[\begin{array}{c}
I_{r} \\
0
\end{array}\right] \Sigma_{1}^{-\frac{1}{2}}=\Sigma_{1}^{-\frac{1}{2}} \Sigma_{1} \Sigma_{1}^{-\frac{1}{2}}=I_{r}
\end{aligned}
$$

$\Longrightarrow V W^{\top}$ is an oblique projector, hence balanced truncation is a Petrov-Galerkin projection method.

## Balanced Truncation

## Properties:

- Reduced-order model is stable with $\mathrm{HSVs} \sigma_{1}, \ldots, \sigma_{r}$.
- Adaptive choice of $r$ via computable error bound:

$$
\|y-\hat{y}\|_{2} \leq\left(2 \sum_{k=r+1}^{n} \sigma_{k}\right)\|u\|_{2} .
$$

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$$

## Balanced Truncation

## Theoretical Background

| Linear, Time-Invariant (LTI) Systems |  |
| :--- | :--- | :--- |
| $\dot{x}=A x+B u$, | $A \in \mathbb{R}^{n \times n}, \quad B \in \mathbb{R}^{n \times m}$, |
| $y=C x+D u, \quad C \in \mathbb{R}^{q \times n}, \quad D \in \mathbb{R}^{q \times m}$. |  |

## Balanced Truncation

## Theoretical Background



## 0000000 <br> Balanced Truncation

## Theoretical Background

```
Linear, Time-Invariant (LTI) Systems
    \(\dot{x}=A x+B u, \quad A \in \mathbb{R}^{n \times n}, \quad B \in \mathbb{R}^{n \times m}\),
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```

State-Space Description for I/O-Relation
Variation-of-constants $\Longrightarrow$

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

## Balanced Truncation

## Theoretical Background

$$
\begin{array}{ll}
\text { Linear, Time-Invariant }(\text { LTI }) \text { Systems } \\
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y=C x+D u, \quad C \in \mathbb{R}^{q \times n}, \quad D \in \mathbb{R}^{q \times m} .
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$$

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


## Balanced Truncation

## Theoretical Background

$$
\begin{array}{lll}
\text { Linear, Time-Invariant }(\text { LTI }) \text { Systems } \\
\dot{x}=A x+B u, & A \in \mathbb{R}^{n \times n}, \quad B \in \mathbb{R}^{n \times m}, \\
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## State-Space Description for I/O-Relation

Variation-of-constants $\Longrightarrow$

$$
\mathcal{S}: u \mapsto y, \quad y(t)=\int_{-\infty}^{t} C e^{A(t-\tau)} B u(\tau) d \tau \quad \text { for all } t \in \mathbb{R} .
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- $\mathcal{S}: \mathcal{U} \rightarrow \mathcal{Y}$ is a linear operator between (function) spaces.
- Recall: $A \in \mathbb{R}^{n \times m}$ is a linear operator, $A: \mathbb{R}^{m} \rightarrow \mathbb{R}^{n}$ !
- Basic Idea: use SVD approximation as for matrix A!
- Problem: in general, $\mathcal{S}$ does not have a discrete SVD and can therefore not be approximated as in the matrix case!


## Balanced Truncation

## Theoretical Background

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But: computationally unfeasible for large-scale systems.

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The Hankel Singular Values are Singular Values!

Theorem
Let $P, Q$ be the controllability and observability Gramians of an LTI system $\Sigma$. Then the Hankel singular values $\Lambda(P Q)^{\frac{1}{2}}=\left\{\sigma_{1}, \ldots, \sigma_{n}\right\}$ are the singular values of the Hankel operator associated to $\Sigma$.

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\mathcal{H}^{*} y_{+}(t)==B^{T} e^{-A^{T} t} \int_{0}^{\infty} e^{A^{T} \tau} C^{T} y_{+}(\tau) d \tau
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Hence,

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z & =\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}} \underbrace{\int_{0}^{\infty} e^{A t} B B^{T} e^{A^{T} t} d t}_{\equiv P} Q z \\
& =\frac{1}{\sigma^{2}} P Q z
\end{aligned}
$$

$$
\Longleftrightarrow \quad P Q z=\sigma^{2} z
$$

$\square$

## Balanced Truncation

The Hankel Singular Values are Singular Values!

## Theorem

Let $P, Q$ be the controllability and observability Gramians of an LTI system $\Sigma$. Then the Hankel singular values $\wedge(P Q)^{\frac{1}{2}}=\left\{\sigma_{1}, \ldots, \sigma_{n}\right\}$ are the singular values of the Hankel operator associated to $\Sigma$.

## Theorem

Let the reduced-order system $\hat{\Sigma}:(\hat{A}, \hat{B}, \hat{C}, \hat{D})$ with $r \leq \hat{n}$ be computed by balanced truncation. Then the reduced-order model $\hat{\Sigma}$ is balanced, stable, minimal, and its HSVs are $\sigma_{1}, \ldots, \sigma_{r}$.

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Proof: Note that in balanced coordinates, the Gramians are diagonal and equal to

$$
\operatorname{diag}\left(\Sigma_{1}, \Sigma_{2}\right)=\operatorname{diag}\left(\sigma_{1}, \ldots, \sigma_{r}, \sigma_{r+1}, \ldots, \sigma_{n}\right)
$$

Hence, the Gramian satisfies
$\left[\begin{array}{ll}A_{11} & A_{12} \\ A_{21} & A_{22}\end{array}\right]\left[\begin{array}{ll}\Sigma_{1} & \\ & \Sigma_{2}\end{array}\right]+\left[\begin{array}{ll}\Sigma_{1} & \\ & \Sigma_{2}\end{array}\right]\left[\begin{array}{ll}A_{11} & A_{12} \\ A_{21} & A_{22}\end{array}\right]^{T}+\left[\begin{array}{l}B_{1} \\ B_{2}\end{array}\right]\left[\begin{array}{l}B_{1} \\ B_{2}\end{array}\right]^{T}=0$,
whence we obtain the "controllability Lyapunov equation" of the reduced-order system,

$$
A_{11} \Sigma_{1}+\Sigma_{1} A_{11}^{T}+B_{1} B_{1}^{T}=0
$$

The result follows from $\hat{A}=A_{11}, \hat{B}=B_{1}, \Sigma_{1}>0$, the solution theory of Lyapunov equations and the analogous considerations for the observability Gramian. (Minimality is a simple consequence of $\hat{P}=\Sigma_{1}=\hat{Q}>0$.)

## Singular Perturbation Approximation (aka Balanced Residualization)

Assume the system

$$
\left[\begin{array}{l}
\dot{x}_{1} \\
\dot{x}_{2}
\end{array}\right]=\left[\begin{array}{ll}
A_{11} & A_{12} \\
A_{21} & A_{22}
\end{array}\right]\left[\begin{array}{l}
x_{1} \\
x_{2}
\end{array}\right]+\left[\begin{array}{l}
B_{1} \\
B_{2}
\end{array}\right] u, \quad y=\left[C_{1}, C_{2}\right]\left[\begin{array}{l}
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Particularly, if $G(0)=\hat{G}(0)$ ("zero steady-state error") is required, one can apply the same condensation technique as in Guyan reduction: instead of $x_{2}=0$, set $\dot{x}_{2}=0$. This yields the reduced-order model

$$
\begin{aligned}
\dot{x}_{1} & =\left(A_{11}-A_{12} A_{22}^{-1} A_{21}\right) x_{1}+\left(B_{1}-A_{12} A_{22}^{-1} B_{2}\right) u \\
y & =\left(C_{1}-C_{2} A_{22}^{-1} A_{21}\right) x_{1}+\left(D-C_{2} A_{22}^{-1} B_{2}\right) u
\end{aligned}
$$

with

- the same properties as the reduced-order model w.r.t. stability, minimality, error bound, but $\hat{D} \neq D$;
- zero steady-state error, $G(0)=\hat{G}(0)$ as desired.


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## Note:

- $A_{22}$ invertible as in balanced coordinates, $A_{22} \Sigma_{2}+\Sigma_{2} A_{22}^{T}+B_{2} B_{2}^{T}=0$ and $\left(A_{22}, B_{2}\right)$ controllable, $\Sigma_{2}>0 \Rightarrow A_{22}$ stable.
- If the original system is not balanced, first compute a minimal realization by applying balanced truncation with $r=\hat{n}$.


## Balancing-Related Methods

## Basic Principle

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

$$
P=Q=\operatorname{diag}\left(\sigma_{1}, \ldots, \sigma_{n}\right)=\Sigma, \quad \sigma_{1} \geq \ldots \geq \sigma_{n}>0
$$ and truncate corresponding realization at size $r$ with $\sigma_{r}>\sigma_{r+1}$.

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## Classical Balanced Truncation (BT)

- $P=$ 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

$$
A P+P A^{T}+B B^{T}=0, \quad A^{T} Q+Q A+C^{T} C=0
$$

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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)

$$
\begin{aligned}
& 0=A P+P A^{T}-P C^{T} C P+B^{T} B \\
& 0=A^{T} Q+Q A-Q B B^{T} Q+C^{T} C .
\end{aligned}
$$

## Balancing-Related Methods

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

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

$$
\hat{A}^{T} Q+Q \hat{A}+Q B_{W}\left(D D^{T}\right)^{-1} B_{W}^{T} Q+C^{T}\left(D D^{T}\right)^{-1} C=0,
$$

where $\hat{A}:=A-B_{W}\left(D D^{T}\right)^{-1} C, B_{W}:=B D^{T}+P C^{T}$.

## Balancing-Related Methods

## Basic Principle

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

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P=Q=\operatorname{diag}\left(\sigma_{1}, \ldots, \sigma_{n}\right)=\Sigma, \quad \sigma_{1} \geq \ldots \geq \sigma_{n}>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

$$
\begin{aligned}
& 0=\bar{A} P+P \bar{A}^{T}+P C^{T} \bar{R}^{-1} C P+B \bar{R}^{-1} B^{T}, \\
& 0=\bar{A}^{T} Q+Q \bar{A}+Q B \bar{R}^{-1} B^{T} Q+C^{T} \bar{R}^{-1} C,
\end{aligned}
$$

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

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## Other Balancing-Based Methods

- Bounded-real balanced truncation (BRBT) - based on bounded real lemma [Opdenacker/Jonckheere '88];
- $H_{\infty}$ balanced truncation (HinfBT) - closed-loop balancing based on $H_{\infty}$ compensator [Mustafa/Glover '91].

Both approaches require solution of dual AREs.

- Frequency-weighted versions of the above approaches.


## Balancing-Related Methods

## Properties

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

$$
\begin{aligned}
\text { BT: }\left\|G-G_{r}\right\|_{\infty} & \leq 2 \sum_{j=r+1}^{n} \sigma_{j}^{B T}, \\
\text { LQGBT: }\left\|G-G_{r}\right\|_{\infty} & \leq 2 \sum_{j=r+1}^{n} \frac{\sigma_{j}^{L Q G}}{\sqrt{1+\left(\sigma_{j}^{L Q G}\right)^{2}}} \\
\text { BST: }\left\|G-G_{r}\right\|_{\infty} & \leq\left(\prod_{j=r+1}^{n} \frac{1+\sigma_{j}^{B S T}}{1-\sigma_{j}^{B S T}}-1\right)\|G\|_{\infty},
\end{aligned}
$$

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


## Outline

(1) Introduction
(2) Mathematical Basics
(3) Model Reduction by ProjectionModal TruncationBalanced Truncation
(6) Solving Large-Scale Matrix Equations

- Linear Matrix Equations
- Numerical Methods for Solving Lyapunov Equations
- Solving Large-Scale Algebraic Riccati Equations
- Software
(7) Final Remarks


## Solving Large-Scale Matrix Equations <br> <br> Large-Scale Algebraic Lyapunov and Riccati Equations

 <br> <br> Large-Scale Algebraic Lyapunov and Riccati Equations}Algebraic Riccati equation (ARE) for $A, G=G^{T}, W=W^{T} \in \mathbb{R}^{n \times n}$ given and $X \in \mathbb{R}^{n \times n}$ unknown:

$$
0=\mathcal{R}(X):=A^{T} X+X A-X G X+W .
$$

$G=0 \Longrightarrow$ Lyapunov equation

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Typical situation in model reduction and optimal control problems for semi-discretized PDEs:

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- $n=10^{3}-10^{6}\left(\Longrightarrow 10^{6}-10^{12}\right.$ unknowns! $)$,
- $A$ has sparse representation $\left(A=-M^{-1} S\right.$ for FEM),
- $G, W$ low-rank with $G, W \in\left\{B B^{\top}, C^{\top} C\right\}$, where $B \in \mathbb{R}^{n \times m}, m \ll n, \quad C \in \mathbb{R}^{p \times n}, p \ll n$.
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## Solving Large-Scale Matrix Equations

## Low-Rank Approximation

Consider spectrum of ARE solution (analogous for Lyapunov equations).
eigenvalues of $\mathrm{P}_{\mathrm{h}}$ for $\mathbf{h}=\mathbf{0 . 0 1}$

## Example:

- Linear 1D heat equation with point control,
- $\Omega=[0,1]$,
- FEM discretization using linear B-splines,
- $h=1 / 100 \Longrightarrow n=101$.


Idea: $X=X^{\top} \geq 0 \Longrightarrow$

$\Longrightarrow$ Goal: compute $Z^{(r)} \in \mathbb{R}^{n \times r}$ directly w/o ever forming $X$ !

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Idea: $X=X^{\top} \geq 0 \Longrightarrow$

$$
X=Z Z^{T}=\sum_{k=1}^{n} \lambda_{k} z_{k} z_{k}^{T} \approx Z^{(r)}\left(Z^{(r)}\right)^{T}=\sum_{k=1}^{r} \lambda_{k} z_{k} z_{k}^{T}
$$

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## Solving Large-Scale Matrix Equations

## Linear Matrix Equations

## Equations without symmetry

Sylvester equation discrete Sylvester equation
$A X+X B=W \quad A X B-X=W$
with data $A \in \mathbb{R}^{n \times n}, B \in \mathbb{R}^{m \times m}, W \in \mathbb{R}^{n \times m}$ and unknown $X \in \mathbb{R}^{n \times m}$.

## Equations with symmetry

Lyapunov equation Stein equation (discrete Lyapunov equation)

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Here: focus on (Sylvester and) Lyapunov equations; analogous results and methods for discrete versions exist.

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## Linear Matrix Equations

## Solvability

Using the Kronecker (tensor) product, $A X+X B=W$ is equivalent to

$$
\left(\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right)\right) \operatorname{vec}(X)=\operatorname{vec}(W)
$$

Hence,

## Sylvester equation has a unique solution

$$
M:=\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right) \text { is invertible. }
$$

$0 \notin \Lambda(M)=\Lambda\left(\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right)\right)=\left\{\lambda_{j}+\mu_{k} \mid \lambda_{j} \in \Lambda(A), \mu_{k} \in \Lambda(B)\right\}$.

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## Corollary

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$$

$$
0 \notin \Lambda(M)=\Lambda\left(\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right)\right)=\left\{\lambda_{j}+\mu_{k} \mid \lambda_{j} \in \Lambda(A), \mu_{k} \in \Lambda(B)\right\} .
$$

$$
\wedge(A) \cap \wedge(-B)=\emptyset
$$

## Corollary

$A, B$ Hurwitz $\Longrightarrow$ Sylvester equation has unique solution.

## Linear Matrix Equations

## Complexity Issues

Solving the Sylvester equation

$$
A X+X B=W
$$

via the equivalent linear system of equations

$$
\left(\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right)\right) \operatorname{vec}(X)=\operatorname{vec}(W)
$$

requires

- LU factorization of $n m \times n m$ matrix; for $n \approx m$, complexity is $\frac{2}{3} n^{6}$;
- storing $n \cdot m$ unknowns: for $n \approx m$ we have $n^{2}$ data for $X$, but up to $n^{4}$ data for triangular factors!

```
Example
n=m=1,000=> Gaussian elimination on an Intel core i7 (Westmere, 6
cores, 3.46 GHz 83.2 GFLOP peak) would take > 94 DAYS and 7.3
TB of memory!
```


## Linear Matrix Equations

## Complexity Issues

Solving the Sylvester equation

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[^3]
## Numerical Methods for Solving Lyapunov Equations Traditional Methods

Bartels-Stewart method for Sylvester and Lyapunov equation (lyap); Hessenberg-Schur method for Sylvester equations (lyap);
Hammarling's method for Lyapunov equations $A X+X A^{T}+G G^{T}=0$ with $A$ Hurwitz (lyapchol).
All based on the fact that if $A, B^{T}$ are in Schur form, then

$$
M=\left(I_{m} \otimes A\right)+\left(B^{T} \otimes I_{n}\right)
$$

is block-upper triangular. Hence, solve $M x=b$ by back-substitution.

- Clever implementation of back-substitution process requires $n m(n+m)$ flops.
- For Sylvester equations, $B$ in Hessenberg form is enough ( $\rightsquigarrow$ Hessenberg-Schur method).
- Hammarling's method computes Cholesky factor $Y$ of $X$ directly.
- All methods require Schur decomposition of $A$ and Schur or Hessenberg decomposition of $B \Rightarrow$ need QR algorithm which requires $25 n^{3}$ flops for Schur decomposition.

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\text { Not feasible for large-scale problems ( } n>10,000 \text { ). }
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## Numerical Methods for Solving Lyapunov Equations The Sign Function Method

## Definition

For $Z \in \mathbb{R}^{n \times n}$ with $\Lambda(Z) \cap \imath \mathbb{R}=\emptyset$ and Jordan canonical form

$$
Z=S\left[\begin{array}{cc}
J^{+} & 0 \\
0 & J^{-}
\end{array}\right] S^{-1}
$$

the matrix sign function is

$$
\operatorname{sign}(Z):=S\left[\begin{array}{cc}
I_{k} & 0 \\
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\end{array}\right] S^{-1}
$$

## Lemma

Let $T \in \mathbb{R}^{n \times n}$ be nonsingular and $Z$ as before, then

$$
\operatorname{sign}\left(T Z T^{-1}\right)=T \operatorname{sign}(Z) T^{-1}
$$

## Numerical Methods for Solving Lyapunov Equations The Sign Function Method

## Computation of sign $(Z)$

$\operatorname{sign}(Z)$ is root of $I_{n} \Longrightarrow$ use Newton's method to compute it:

$$
\begin{aligned}
& Z_{0} \leftarrow Z, \quad Z_{j+1} \leftarrow \frac{1}{2}\left(c_{j} Z_{j}+\frac{1}{c_{j}} Z_{j}^{-1}\right), \quad j=1,2, \ldots \\
\Longrightarrow & \operatorname{sign}(Z)=\lim _{j \rightarrow \infty} Z_{j} .
\end{aligned}
$$

$c_{j}>0$ is scaling parameter for convergence acceleration and rounding error minimization, e.g.

$$
c_{j}=\sqrt{\frac{\left\|Z_{j}^{-1}\right\|_{F}}{\left\|Z_{j}\right\|_{F}}}
$$

based on "equilibrating" the norms of the two summands [Higham '86].

## Solving Lyapunov Equations with the Matrix Sign Function Method

Key observation:
If $X \in \mathbb{R}^{n \times n}$ is a solution of $A X+X A^{T}+W=0$, then

$$
\underbrace{\left[\begin{array}{cc}
I_{n} & -X \\
0 & I_{n}
\end{array}\right]}_{=T^{-1}} \underbrace{\left[\begin{array}{cc}
A & W \\
0 & -A^{T}
\end{array}\right]}_{=: H} \underbrace{\left[\begin{array}{cc}
I_{n} & X \\
0 & I_{n}
\end{array}\right]}_{=: T}=\left[\begin{array}{cc}
A & 0 \\
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$$

Hence, if $A$ is Hurwitz (i.e., asymptotically stable), then


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0 & -A^{T}
\end{array}\right] .
$$

Hence, if $A$ is Hurwitz (i.e., asymptotically stable), then

$$
\begin{aligned}
\operatorname{sign}(H) & =\operatorname{sign}\left(T\left[\begin{array}{cc}
A & 0 \\
0 & -A^{T}
\end{array}\right] T^{-1}\right)=T \operatorname{sign}\left(\left[\begin{array}{cc}
A & 0 \\
0 & -A^{T}
\end{array}\right]\right) T^{-1} \\
& =\left[\begin{array}{cc}
-I_{n} & 2 X \\
0 & I_{n}
\end{array}\right]
\end{aligned}
$$

## Solving Lyapunov Equations with the Matrix Sign Function Method

Apply sign function iteration $Z \leftarrow \frac{1}{2}\left(Z+Z^{-1}\right)$ to $H=\left[\begin{array}{cc}A & W \\ 0 & -A^{T}\end{array}\right]$ :

$$
H+H^{-1}=\left[\begin{array}{cc}
A & W \\
0 & -A^{T}
\end{array}\right]+\left[\begin{array}{cc}
A^{-1} & A^{-1} W A^{-T} \\
0 & -A^{-T}
\end{array}\right]
$$

$\Longrightarrow$ Sign function iteration for Lyapunov equation:

$$
\begin{array}{ll}
A_{0} \leftarrow A, & A_{j+1} \leftarrow \frac{1}{2}\left(A_{j}+A_{j}^{-1}\right), \\
W_{0} \leftarrow G, & W_{i+1} \leftarrow \frac{1}{2}\left(W_{i}+A_{i}^{-1} W_{i} A_{i}^{-T}\right),
\end{array} \quad j=0,1,2, \ldots
$$

Define $A_{\infty}:=\lim _{j \rightarrow \infty} A_{j}, W_{\infty}:=\lim _{j \rightarrow \infty} W_{j}$.

## Theorem

If $A$ is Hurwitz, then

$$
A_{\infty}=-I_{n} \quad \text { and } \quad X=\frac{1}{2} W_{\infty} .
$$

## Solving Lyapunov Equations with the Matrix Sign Function Method

 Factored formRecall sign function iteration for $A X+X A^{T}+W=0$ :

$$
\begin{array}{ll}
A_{0} \leftarrow A, & A_{j+1} \leftarrow \frac{1}{2}\left(A_{j}+A_{j}^{-1}\right), \\
W_{0} \leftarrow G, & W_{j+1} \leftarrow \frac{1}{2}\left(W_{j}+A_{j}^{-1} W_{j} A_{j}^{-T}\right),
\end{array} \quad j=0,1,2, \ldots
$$

Now consider the second iteration for $W=B B^{\top}$, starting with $W_{0}=B B^{T}=: B_{0} B_{0}^{T}:$

$$
\begin{aligned}
\frac{1}{2}\left(W_{j}+A_{j}^{-1} W_{j} A_{j}^{-\top}\right) & =\frac{1}{2}\left(B_{j} B_{j}^{\top}+A_{j}^{-1} B_{j} B_{j}^{\top} A_{j}^{-\top}\right) \\
& =\frac{1}{2}\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right]\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right]^{\top}
\end{aligned}
$$

Hence, obtain factored iteration

$$
B_{j+1} \leftarrow \frac{1}{\sqrt{2}}\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right]
$$

with $S:=\frac{1}{\sqrt{2}} \lim _{j \rightarrow \infty} B_{j}$ and $X=S S^{\top}$

## Solving Lyapunov Equations with the Matrix Sign Function Method

 Factored formRecall sign function iteration for $A X+X A^{T}+W=0$ :

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\begin{array}{ll}
A_{0} \leftarrow A, & A_{j+1} \leftarrow \frac{1}{2}\left(A_{j}+A_{j}^{-1}\right), \\
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Solving Lyapunov Equations with the Matrix Sign Function Method Factored form

Recall sign function iteration for $A X+X A^{T}+W=0$ :

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A_{0} \leftarrow A, & A_{j+1} \leftarrow \frac{1}{2}\left(A_{j}+A_{j}^{-1}\right), \\
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$$

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$$
\begin{aligned}
\frac{1}{2}\left(W_{j}+A_{j}^{-1} W_{j} A_{j}^{-T}\right) & =\frac{1}{2}\left(B_{j} B_{j}^{T}+A_{j}^{-1} B_{j} B_{j}^{T} A_{j}^{-T}\right) \\
& =\frac{1}{2}\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right]\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right]^{T}
\end{aligned}
$$

Hence, obtain factored iteration

with $S:=\frac{1}{\sqrt{2}} \lim _{j \rightarrow \infty} B_{j}$ and $X=S S^{T}$

## Solving Lyapunov Equations with the Matrix Sign Function Method Factored form

Recall sign function iteration for $A X+X A^{T}+W=0$ :

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\end{array}
$$

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with $S:=\frac{1}{\sqrt{2}} \lim _{j \rightarrow \infty} B_{j}$ and $X=S S^{T}$.

## Solving Lyapunov Equations with the Matrix Sign Function Method Factored form <br> [B./Quintana-Ortí '97]

Factored sign function iteration for $A\left(S S^{T}\right)+\left(S S^{T}\right) A^{T}+B B^{T}=0$

$$
\begin{array}{ll}
A_{0} \leftarrow A, & A_{j+1} \leftarrow \frac{1}{2}\left(A_{j}+A_{j}^{-1}\right), \\
B_{0} \leftarrow B, & B_{j+1} \leftarrow \frac{1}{\sqrt{2}}\left[\begin{array}{ll}
B_{j} & A_{j}^{-1} B_{j}
\end{array}\right],
\end{array} \quad j=0,1,2, \ldots .
$$

## Remarks:

- To get both Gramians, run in parallel

$$
C_{j+1} \leftarrow \frac{1}{\sqrt{2}}\left[\begin{array}{c}
C_{j} \\
C_{j} A_{j}^{-1}
\end{array}\right]
$$

- To avoid growth in numbers of columns of $B_{j}$ (or rows of $C_{j}$ ): column compression by RRLQ or truncated SVD.
- Several options to incorporate scaling, e.g., scale " $A$ "-iteration only.
- Simple stopping criterion: $\left\|A_{j}+I_{n}\right\|_{F} \leq$ tol.


# 0000000000000000000000000000000 

## Numerical Methods for Solving Lyapunov Equations The ADI Method

Recall Peaceman Rachford ADI:
Consider $A u=s$ 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

$$
\begin{aligned}
& (H+p l) v=r \\
& (V+p l) w=t
\end{aligned}
$$

can be solved easily/efficiently.

## AD Iteration

If $H, V$ spd $\Rightarrow \exists p_{k}, k=1,2, \ldots$ such that

$$
\begin{aligned}
u_{0} & =0 \\
\left(H+p_{k} /\right) u_{k-\frac{1}{2}} & =\left(p_{k} l-V\right) u_{k-1}+s \\
\left(V+p_{k} I\right) u_{k} & =\left(p_{k} l-H\right) u_{k-\frac{1}{2}}+s
\end{aligned}
$$

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

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## Numerical Methods for Solving Lyapunov Equations

The Lyapunov operator

$$
\mathcal{L}: \quad X \quad \mapsto \quad A X+X A^{T}
$$

can be decomposed into the linear operators

$$
\mathcal{L}_{H}: X \mapsto A X, \quad \mathcal{L}_{V}: X \mapsto X A^{T} .
$$

In analogy to the standard ADI method we find the

## ADI iteration for the Lyapunov equation

$$
\begin{aligned}
X_{0} & =0 \\
\left(A+p_{k} I\right) X_{k-\frac{1}{1}} & =-W-X_{k-1}\left(A^{T}-p_{k} I\right) \\
\left(A+p_{k} I\right) X_{k}^{T} & =-W-X_{k-\frac{1}{2}}^{T}\left(A^{T}-p_{k} I\right) .
\end{aligned}
$$

# 0000000000000000000000000000000 

## Numerical Methods for Solving Lyapunov Equations Low-Rank ADI

Consider $A X+X A^{T}=-B B^{T}$ for stable $A ; B \in \mathbb{R}^{n \times m}$ with $m \ll n$.

## ADI iteration for the Lyapunov equation

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

$$
\begin{array}{ccc}
X_{0} & = & 0 \\
\left(A+p_{k} I\right) X_{k-\frac{1}{2}} & = & -B B^{T}-X_{k-1}\left(A^{T}-p_{k} I\right) \\
\left(A+p_{k} I\right) X_{k}^{T^{2}} & = & -B B^{T}-X_{k-\frac{1}{2}}^{T}\left(A^{T}-p_{k} I\right)
\end{array}
$$

Rewrite as one step iteration and factorize $X_{k}=Z_{k} Z_{k}^{\top}, k=0, \ldots, k_{\max }$

$$
\begin{aligned}
Z_{0} Z_{0}^{T}= & 0 \\
Z_{k} Z_{k}^{\top}= & -2 p_{k}\left(A+p_{k} I\right)^{-1} B B^{T}\left(A+p_{k} I\right)^{-T} \\
& +\left(A+p_{k} I\right)^{-1}\left(A-p_{k} I\right) Z_{k-1} Z_{k-1}^{T}\left(A-p_{k} I\right)^{T}\left(A+p_{k} I\right)^{-T}
\end{aligned}
$$

$\rightsquigarrow$ 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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& +\left(A+p_{k} I\right)^{-1}\left(A-p_{k} I\right) Z_{k-1} Z_{k-1}^{T}\left(A-p_{k} I\right)^{T}\left(A+p_{k} I\right)^{-T}
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\left(A+p_{k} I\right) X_{k}^{T^{2}} & = & -B B^{T}-X_{k-\frac{1}{2}}^{T}\left(A^{T}-p_{k} I\right)
\end{array}
$$

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$$
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& +\left(A+p_{k} I\right)^{-1}\left(A-p_{k} I\right) Z_{k-1} Z_{k-1}^{T}\left(A-p_{k} I\right)^{T}\left(A+p_{k} I\right)^{-T}
\end{aligned}
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[Penzl '97/'00, Li/White '99/'02, B./Li/Penzl '99/'08, Gugercin/Sorensen/Antoulas '03]

# 0,000000000000000000000000000000 

## Solving Large-Scale Matrix Equations

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

[Penzl '00]
Observing that $\left(A-p_{i} l\right),\left(A+p_{k} l\right)^{-1}$ commute, we rewrite $Z_{k_{\max }}$ as
$Z_{k_{\max }}=\left[z_{k_{\max }}, P_{k_{\max }-1} z_{k_{\max }}, P_{k_{\max }-2}\left(P_{k_{\max }-1} z_{k_{\max }}\right), \ldots, P_{1}\left(P_{2} \ldots P_{k_{\max }-1} z_{k_{\max }}\right)\right]$,
where

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

and

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

## Solving Large-Scale Matrix Equations Numerical Methods for Solving Lyapunov Equations

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

[Penzl '00]
Observing that $\left(A-p_{i} I\right),\left(A+p_{k} I\right)^{-1}$ commute, we rewrite $Z_{k_{\max }}$ as

$$
Z_{k_{\max }}=\left[z_{k_{\max }}, P_{k_{\max }-1} z_{k_{\max }}, P_{k_{\max }-2}\left(P_{k_{\max }-1} z_{k_{\max }}\right), \ldots, P_{1}\left(P_{2} \ldots P_{k_{\max }-1} z_{k_{\max }}\right)\right]
$$

[LI/White '02]
where

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

and

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

## Numerical Methods for Solving Lyapunov Equations

 Lyapunov equation $0=A X+X A^{T}+B B^{T}$.Algorithm [Penzl '97/'00, Li/White '99/'02, B. 04, B./Li/Penzl '99/'08]

$$
\begin{aligned}
& V_{1} \leftarrow \sqrt{-2 \text { re } p_{1}}\left(A+p_{1} /\right)^{-1} B, \quad Z_{1} \leftarrow V_{1} \\
& \text { FOR } k=2,3, \ldots
\end{aligned}
$$

$$
\begin{aligned}
& V_{k} \leftarrow \sqrt{\frac{r e}{r e} p_{k}}\left(V_{k-1}-\left(p_{k}+\overline{p_{k-1}}\right)\left(A+p_{k} I\right)^{-1} V_{k-1}\right) \\
& Z_{k} \leftarrow\left[Z_{k-1} \quad V_{k}\right] \\
& Z_{k} \leftarrow \operatorname{rrlq}\left(Z_{k}, \tau\right) \quad \text { column compression }
\end{aligned}
$$

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


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

$$
\left.z_{k_{\max }}=\left[\begin{array}{lll}
v_{1} & \ldots & v_{k_{\max }}
\end{array}\right], \quad v_{k}=\right] \in \mathbb{C}^{n \times m} .
$$

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].

## Numerical Results for ADI

## Optimal Cooling of Steel Profiles

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

$$
\begin{aligned}
c \cdot \rho \frac{\partial}{\partial t} x & =\lambda \Delta x, \quad \xi \in \Omega \\
\lambda \frac{\partial}{\partial n} x & =\kappa\left(u_{k}-x\right), \quad \xi \in \Gamma_{k}, 1 \leq k \leq 7 \\
\frac{\partial}{\partial n} x & =0, \quad \xi \in \Gamma_{7} . \\
\Longrightarrow m=7, q & =6 .
\end{aligned}
$$

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


Source: Physical model: courtesy of Mannesmann/Demag.
Math. model: Tröltzsch/Unger 1999/2001, Penzl 1999, SaAk 2003.

## Numerical Results for ADI

## Optimal Cooling of Steel Profiles

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

$$
A P M^{T}+M P A^{T}+B B^{T}=0, \quad A^{T} Q M+M^{T} Q A+C^{T} C=0
$$

for $n=79,841$.

- 25 shifts chosen by Penzl heuristic from 50/25 Ritz values of $A$ of largest/smallest magnitude, no column compression performed.
- No factorization of mass matrix required.
- Computations done on Core2Duo at 2.8 GHz with 3GB RAM and 32Bit-MATLAB.



CPU times: $626 / 356 \mathrm{sec}$.

## Numerical Results for ADI

Scaling

Computations by Martin Köhler '10

- $A \in \mathbb{R}^{n \times n} \equiv$ FDM matrix for 2D heat equation on $[0,1]^{2}$ (LyAPACK benchmark demo_11, $m=1$ ).
- 16 shifts chosen by Penzl heuristic from 50/25 Ritz values of $A$ of largest/smallest magnitude.
- Computations on 2 dual core Intel Xeon 5160 with 16 GB RAM using M.E.S.S. (http://svncsc.mpi-magdeburg.mpg.de/trac/messtrac/).


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CPU Times

| n | M.E.S.S. ${ }^{1}(\mathrm{C})$ | 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 |

## Numerical Results for ADI

Scaling

Computations by Martin Köhler '10

- $A \in \mathbb{R}^{n \times n} \equiv$ 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 on 2 dual core Intel Xeon 5160 with 16 GB RAM using M.E.S.S. (http://svncsc.mpi-magdeburg.mpg.de/trac/messtrac/).


Note: for $n=1,000,000$, first sparse LU needs $\sim 1,100$ sec., using UMFPACK this reduces to 30 sec .

## Factored Galerkin-ADI Iteration

## Lyapunov equation $0=A X+X A^{T}+B B^{T}$

Projection-based methods for Lyapunov equations with $A+A^{T}<0$ :
© Compute orthonormal basis range $(Z), Z \in \mathbb{R}^{n \times r}$, for subspace $\mathcal{Z} \subset \mathbb{R}^{n}$, $\operatorname{dim} \mathcal{Z}=r$.
(2) Set $\hat{A}:=Z^{\top} A Z, \hat{B}:=Z^{\top} B$.
(0) 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 $m=1$ :

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

[Sadd '90, Jaimoukha/Kasenally '94, Jbilou '02-'08].

- K-PIK [Simoncini $\left.{ }^{\circ} 07\right]$,

$$
\mathcal{Z}=\mathcal{K}(A, B, r) \cup \mathcal{K}\left(A^{-1}, B, r\right) .
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- Rational Krylov [Druskin/Simoncini '11].


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(9) Use $X \approx Z \hat{X} Z^{T}$.

## Examples:

- ADI subspace [B./R.-C. Li/Truhar '08]:

$$
\mathcal{Z}=\operatorname{colspan}\left[\begin{array}{lll}
V_{1}, & \ldots, & V_{r}
\end{array}\right]
$$

Note:
(1) ADI subspace is rational Krylov subspace [J.-R. Li/White '02].
(2) Similar approach: ADI-preconditioned global Arnoldi method [Jbilou '08].

## Numerical Methods for Solving Lyapunov Equations Numerical examples for Galerkin-ADI

FEM semi-discretized control problem for parabolic PDE:

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


## Good ADI shifts



CPU times: 80s (projection every 5th ADI step) vs. 94s (no projection).
Computations by Jens Saak '10.

## Numerical Methods for Solving Lyapunov Equations Numerical examples for Galerkin-ADI

FEM semi-discretized control problem for parabolic PDE:

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


## Bad ADI shifts



CPU times: 368s (projection every 5th ADI step) vs. 1207s (no projection).
Computations by Jens Saak '10.

## Numerical Methods for Solving Lyapunov Equations

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

## M.E.S.S. w/o Galerkin projection and column compression




Rank of solution factors: 532 / 426

## M.E.S.S. with Galerkin projection and column compression




Rank of solution factors: 269 / 205

## Solving Large-Scale Matrix Equations

Numerical example for BT: Optimal Cooling of Steel Profiles

## $n=1,357$, Absolute Error



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


## Solving Large-Scale Matrix Equations

## Numerical example for BT: Optimal Cooling of Steel Profiles

## $n=1,357$, Absolute Error



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


## $n=79,841$, Absolute Error



- BT model computed using M.E.S.S. in MATLAB,
- dualcore, computation time: $<10 \mathrm{~min}$.


## Solving Large-Scale Matrix Equations

## 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.imtek.de/simulation/benchmark
Courtesy of D. Billger (Imego Institute, Göteborg), Saab Bofors Dynamics AB.

## Solving Large-Scale Matrix Equations

Numerical example for BT: Microgyroscope (Butterfly Gyro)

- FEM discretization of structure dynamical model using quadratic tetrahedral elements (ANSYS-SOLID187)
$\rightsquigarrow n=34,722, m=1, q=12$.
- Reduced model computed using SpaRed, $r=30$.


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## Frequency Repsonse Analysis



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## Frequency Repsonse Analysis



## Hankel Singular Values



## Solving Large-Scale Algebraic Riccati Equations

## Theorem

Consider the (continuous-time) algebraic Riccati equation (ARE)

$$
0=\mathcal{R}(X)=C^{T} C+A^{T} X+X A-X B B^{T} X
$$

with $A \in \mathbb{R}^{n \times n}, B \in \mathbb{R}^{n \times m}, C \in \mathbb{R}^{q \times n},(A, B)$ stabilizable, $(A, C)$ detectable. Then:
(a) There exists a unique stabilizing $X_{*} \in\left\{X \in \mathbb{R}^{n \times n} \mid \mathcal{R}(X)=0\right\}$, i.e., $\Lambda\left(A-B B^{T} X_{*}\right) \in \mathbb{C}^{-}$.
(b) $X_{*}=X_{*}^{\top} \geq 0$ and $X_{*} \geq X$ for all $X \in\left\{X \in \mathbb{R}^{n \times n} \mid \mathcal{R}(X)=0\right\}$.
(c) If $(A, C)$ observable, then $X_{*}>0$.
(d) $\operatorname{span}\left\{\left[\begin{array}{c}I_{n} \\ -X_{*}\end{array}\right]\right\}$ is the unique invariant subspace of the Hamiltonian matrix

$$
H=\left[\begin{array}{cc}
A & B B^{T} \\
C^{T} C & -A^{T}
\end{array}\right]
$$

corresponding to $\Lambda(H) \cap \mathbb{C}^{-}$.

## Solving Large-Scale Algebraic Riccati Equations

## Numerical Methods (incomplete list)

- Invariant subspace methods ( $\rightsquigarrow$ eigenproblem for Hamiltonian matrix):
- Schur vector method (care)
- Hamiltonian SR algorithm
- Symplectic URV-based method

> [B./Mehrmann/Xu '97/'98, Chu/Liu/Mehrmann '07]

- Spectral projection methods
- Sign function method
- Disk function method
[Roberts '71, Byers '87] [Bai/Demmel/Gu '94, B. '97]
- (rational, global) Krylov subspace techniques
[Jaimoukha/Kasenally '94, Jbilou '03/'06, Heyouni/Jbilou '09]
- Newton's method
- Kleinman iteration
[Kleinman '68]
- Line search acceleration
[B./Byers '98]
- Newton-ADI
[B./J.-R. Li/Penzl '99/'08]
- Inexact Newton
[Feitzinger/Hylla/Sachs '09,B./Heinkenschloss/SaAK/Weichelt '15]


## Solving Large-Scale Matrix Equations

## Software

## Lyapack

MATLAB toolbox for solving

- Lyapunov equations and algebraic Riccati equations,
- model reduction and LQR problems.

Main work horse: Low-rank ADI and Newton-ADI iterations.

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## M.E.S.S. - Matrix Equations Sparse Solvers

- Extended and revised version of Lyapack.
- Includes solvers for large-scale differential Riccati equations (based on Rosenbrock and BDF methods).
- Many algorithmic improvements:
- new ADI parameter selection,
- column compression based on RRQR,
- more efficient use of direct solvers,
- treatment of generalized systems without factorization of the mass matrix,
- new ADI versions avoiding complex arithmetic etc.
- $C$ and MATLAB versions.


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## Topics Not Covered

- Special methods for second-order (mechanical) systems.
- Extensions to bilinear and stochastic systems.
- Balanced truncation for discrete-time systems.
- Extensions to descriptor systems $E \dot{x}=A x+B u, E$ singular.
- Frequency-limited/-weighted balanced truncation.
- Application to 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.

## Further Reading - Balanced Truncation

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Computational methods for linear matrix equations (survey article).
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http://www.dm.unibo.it/-simoncin/matrixeq.pdf.


[^0]:    Source: http://en.wikipedia.org/wiki/File:Transistor_Count_and_Moore'sLaw_-_2011.svg

[^1]:    Corollary
    $A, B$ Hurwitz $=\Longrightarrow$ Sylvester equation has unique solution.

[^2]:    Corollary
    $A, B$ Hurwitz $\Longrightarrow$ Sylvester equation has unique solution.

[^3]:    Example
    $n=m=1,000 \Rightarrow$ Gaussian elimination on an Intel core i7 (Westmere, 6 cores, $3.46 \mathrm{GHz} \rightsquigarrow 83.2$ GFLOP peak) would take $>94$ DAYS and 7.3 TB of memory!

